

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081623\
 Data File : BM041422.D
 Acq On : 16 Aug 2023 13:25
 Operator : MA/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 08/17/2023
 Supervised By :mohammad ahmed 08/17/2023

Quant Time: Aug 16 16:50:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 16:44:10 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	165857	20.000	ng	0.00
21) Naphthalene-d8	10.569	136	650935	20.000	ng	0.00
39) Acenaphthene-d10	14.427	164	415557	20.000	ng	0.00
64) Phenanthrene-d10	17.186	188	893635	20.000	ng	0.00
76) Chrysene-d12	21.392	240	727190	20.000	ng	0.00
86) Perylene-d12	23.750	264	690539	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	780170	78.252	ng	0.00
7) Phenol-d6	6.951	99	1056281	78.024	ng	0.00
23) Nitrobenzene-d5	8.945	82	1032503	75.718	ng	0.00
42) 2,4,6-Tribromophenol	15.933	330	494938	75.051	ng	0.00
45) 2-Fluorobiphenyl	13.045	172	2382943	76.841	ng	0.00
79) Terphenyl-d14	19.827	244	3556914	85.117	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.246	88	159893	37.748	ng	100
3) Pyridine	3.652	79	439401m	40.266	ng	
4) n-Nitrosodimethylamine	3.569	42	198066	38.366	ng	100
6) Aniline	7.104	93	634884	39.437	ng	100
8) 2-Chlorophenol	7.328	128	395264	38.373	ng	100
9) Benzaldehyde	6.916	77	270805	35.667	ng	100
10) Phenol	6.975	94	511899	39.136	ng	100
11) bis(2-Chloroethyl)ether	7.204	93	399212	38.152	ng	100
12) 1,3-Dichlorobenzene	7.651	146	436709	38.146	ng	100
13) 1,4-Dichlorobenzene	7.798	146	444641	38.414	ng	100
14) 1,2-Dichlorobenzene	8.116	146	431475	38.620	ng	100
15) Benzyl Alcohol	8.028	79	390587	40.640	ng	100
16) 2,2'-oxybis(1-Chloropr...	8.304	45	469720	37.462	ng	100
17) 2-Methylphenol	8.222	107	361532	38.666	ng	100
18) Hexachloroethane	8.828	117	166648	37.914	ng	100
19) n-Nitroso-di-n-propyla...	8.587	70	354858	38.208	ng	100
20) 3+4-Methylphenols	8.557	107	493801	38.452	ng	100
22) Acetophenone	8.604	105	626010	37.319	ng	100
24) Nitrobenzene	8.986	77	519569	38.158	ng	100
25) Isophorone	9.510	82	976882	39.139	ng	100
26) 2-Nitrophenol	9.692	139	234587	39.459	ng	100
27) 2,4-Dimethylphenol	9.757	122	374725	39.145	ng	100
28) bis(2-Chloroethoxy)met...	9.992	93	557132	38.349	ng	100
29) 2,4-Dichlorophenol	10.228	162	406860	38.796	ng	100
30) 1,2,4-Trichlorobenzene	10.433	180	442306	37.984	ng	100
31) Naphthalene	10.616	128	1297279	38.331	ng	100
32) Benzoic acid	9.951	122	308754	40.158	ng	100
33) 4-Chloroaniline	10.751	127	562764	39.387	ng	100
34) Hexachlorobutadiene	10.886	225	290974	37.785	ng	100
35) Caprolactam	11.580	113	127710	39.282	ng	100
36) 4-Chloro-3-methylphenol	11.886	107	435897	39.055	ng	100
37) 2-Methylnaphthalene	12.239	142	946276	38.426	ng	100
38) 1-Methylnaphthalene	12.457	142	899124	38.646	ng	100
40) 1,2,4,5-Tetrachloroben...	12.604	216	519089	38.287	ng	100
41) Hexachlorocyclopentadiene	12.569	237	250722	37.671	ng	100
43) 2,4,6-Trichlorophenol	12.863	196	352707	38.872	ng	100

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081623\
 Data File : BM041422.D
 Acq On : 16 Aug 2023 13:25
 Operator : MA/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 08/17/2023
 Supervised By :mohammad ahmed 08/17/2023

Quant Time: Aug 16 16:50:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 16:44:10 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.939	196	407062	38.753	ng	100
46) 1,1'-Biphenyl	13.257	154	1212131	38.650	ng	100
47) 2-Chloronaphthalene	13.304	162	927521	38.480	ng	100
48) 2-Nitroaniline	13.527	65	320767	40.727	ng	100
49) Acenaphthylene	14.151	152	1515097	38.789	ng	100
50) Dimethylphthalate	13.904	163	1239163	38.188	ng	100
51) 2,6-Dinitrotoluene	14.033	165	271360	39.021	ng	100
52) Acenaphthene	14.492	154	898595	38.465	ng	100
53) 3-Nitroaniline	14.368	138	276506	39.869	ng	100
54) 2,4-Dinitrophenol	14.580	184	170423	39.180	ng	100
55) Dibenzofuran	14.833	168	1494821	38.158	ng	100
56) 4-Nitrophenol	14.686	139	206639	39.568	ng	100
57) 2,4-Dinitrotoluene	14.827	165	368964	38.822	ng	100
58) Fluorene	15.480	166	1190044	37.906	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.063	232	343311	38.457	ng	100
60) Diethylphthalate	15.263	149	1244333	38.381	ng	100
61) 4-Chlorophenyl-phenyle...	15.480	204	637362	37.656	ng	100
62) 4-Nitroaniline	15.539	138	257105	40.637	ng	100
63) Azobenzene	15.774	77	1289629	39.480	ng	100
65) 4,6-Dinitro-2-methylph...	15.592	198	231257	40.426	ng	100
66) n-Nitrosodiphenylamine	15.704	169	1022489	38.870	ng	100
67) 4-Bromophenyl-phenylether	16.374	248	394526	38.526	ng	100
68) Hexachlorobenzene	16.486	284	426393	38.154	ng	100
69) Atrazine	16.662	200	291525	39.053	ng	100
70) Pentachlorophenol	16.839	266	307226	39.259	ng	100
71) Phenanthrene	17.233	178	1818450	38.549	ng	100
72) Anthracene	17.321	178	1856979	38.821	ng	100
73) Carbazole	17.604	167	1656984	38.932	ng	100
74) Di-n-butylphthalate	18.162	149	2080883	38.918	ng	100
75) Fluoranthene	19.256	202	2133819	37.994	ng	100
77) Benzidine	19.462	184	353851	42.451	ng	100
78) Pyrene	19.621	202	2165091	41.631	ng	100
80) Butylbenzylphthalate	20.527	149	852791	40.940	ng	100
81) Benzo(a)anthracene	21.380	228	1912382	38.510	ng	100
82) 3,3'-Dichlorobenzidine	21.321	252	628119	39.108	ng	100
83) Chrysene	21.433	228	1797128	38.072	ng	100
84) Bis(2-ethylhexyl)phtha...	21.297	149	1159691	39.188	ng	100
85) Di-n-octyl phthalate	22.215	149	1781991	36.827	ng	100
87) Indeno(1,2,3-cd)pyrene	26.197	276	1707487	37.978	ng	100
88) Benzo(b)fluoranthene	23.039	252	1629068	39.592	ng	100
89) Benzo(k)fluoranthene	23.086	252	1583253	38.065	ng	100
90) Benzo(a)pyrene	23.650	252	1506918	38.451	ng	100
91) Dibenzo(a,h)anthracene	26.203	278	1425565	37.701	ng	100
92) Benzo(g,h,i)perylene	26.944	276	1412073	38.630	ng	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Manual Integrations

Reviewed By :Yogesh Patel 08/17/2023
Supervised By :mohammad ahmed 08/17/2023

Abundance

TIC: BM041422.D\data.ms

Time-->

1,4-Dioxane
n-Nitrosodimethylamine,
2-Fluorophenol,S
Benzaldehyde,C
Phenol-d6,S
Bis(2-chlorophenyl)ether
1,4-Dichlorobenzene,C
Benzaldehyde,C
2,2'-oxybis(1-chloroethanol)
3,4-Methylenediphenylamine,P
Hexachlorocyclopentadiene,Nitrobenzene-d5,S
Isophorone
2-Nitrophenol
Benzoic acid,bis(2-chloroethoxy)methane
2,4-Dichlorobenzene
Naphthalene
Hexachlorobutadiene,C
Caprolactam
4-Chloro-3-methylphenol,C
2-Methylnaphthalene
Hexachlorocyclopentadiene,Nitrobenzene
2,4,5-Trichlorophenol,C
2-Chloronaphthalene
2-Nitroaniline
2,6-Dinitrophenol
Acenaphthylene
Acenaphthene,C
3-Nitrophenol
4-Nitrophenol
2,3,4,6-Tetrachlorophenol
Diethylphthalate
Dibenzofuran
Azobenzene,Nitrosodiphenylamine,c
2,4,6-Trichlorophenol,S
Hexachlorocyclopentadiene,phenylether
Alkylchlorophenol,C
Phenanthrene-d10,L
Carbazole
Di-n-butylphthalate
Benzidine
Fluoranthene,C
Pyrene
Butylbenzylphthalate
Bis(2-ethylhexyl)phthalate
Bis(2-ethylhexyl)terephthalate
Di-n-octyl phthalate,c
Benz(a)fluoranthene
Benzo(a)pyrene,C
Perylene-d12,L
Dibenz(a,h)pyrene
Benzo(g,h,i)perylene