

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072823\
 Data File : BM041160.D
 Acq On : 28 Jul 2023 11:29
 Operator : MA/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/29/2023
 Supervised By :mohammad ahmed 07/29/2023

Quant Time: Jul 29 00:03:52 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 28 23:54:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	156440	20.000	ng	0.00
21) Naphthalene-d8	10.610	136	591141	20.000	ng	0.00
39) Acenaphthene-d10	14.468	164	321577	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	647514	20.000	ng	0.00
76) Chrysene-d12	21.427	240	616295	20.000	ng	0.00
86) Perylene-d12	23.803	264	703039	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	217719	20.800	ng	0.00
7) Phenol-d6	6.981	99	275742	20.323	ng	0.00
23) Nitrobenzene-d5	8.981	82	248072	19.516	ng	0.00
42) 2,4,6-Tribromophenol	15.968	330	82848	18.214	ng	0.00
45) 2-Fluorobiphenyl	13.086	172	496257	19.494	ng	0.00
79) Terphenyl-d14	19.862	244	633361	18.581	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.269	88	47512	10.611	ng	97
3) Pyridine	3.681	79	117064	9.974	ng	98
4) n-Nitrosodimethylamine	3.593	42	55725	10.234	ng	# 96
6) Aniline	7.140	93	160334	10.098	ng	99
8) 2-Chlorophenol	7.369	128	105184	10.294	ng	99
9) Benzaldehyde	6.957	77	84915	11.848	ng	97
10) Phenol	7.004	94	134331	10.251	ng	99
11) bis(2-Chloroethyl)ether	7.239	93	110737	10.439	ng	99
12) 1,3-Dichlorobenzene	7.692	146	116903	10.510	ng	99
13) 1,4-Dichlorobenzene	7.839	146	117423	10.503	ng	100
14) 1,2-Dichlorobenzene	8.157	146	112955	10.537	ng	98
15) Benzyl Alcohol	8.063	79	78354	9.145	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.345	45	146753m	10.384	ng	
17) 2-Methylphenol	8.263	107	88887	9.928	ng	99
18) Hexachloroethane	8.875	117	41543	9.981	ng	97
19) n-Nitroso-di-n-propyla...	8.622	70	81112	9.842	ng	99
20) 3+4-Methylphenols	8.592	107	119191	9.957	ng	96
22) Acetophenone	8.645	105	151762	9.694	ng	# 98
24) Nitrobenzene	9.022	77	126983	9.878	ng	95
25) Isophorone	9.545	82	201531	9.340	ng	100
26) 2-Nitrophenol	9.733	139	43749	8.706	ng	99
27) 2,4-Dimethylphenol	9.798	122	85842	9.708	ng	99
28) bis(2-Chloroethoxy)met...	10.033	93	132375	9.794	ng	99
29) 2,4-Dichlorophenol	10.269	162	84377	9.462	ng	96
30) 1,2,4-Trichlorobenzene	10.475	180	96856	9.831	ng	96
31) Naphthalene	10.663	128	321358	10.018	ng	99
32) Benzoic acid	9.898	122	34969	11.636	ng	97
33) 4-Chloroaniline	10.798	127	123406	9.626	ng	99
34) Hexachlorobutadiene	10.933	225	56834	9.591	ng	96
35) Caprolactam	11.580	113	21055m	8.162	ng	
36) 4-Chloro-3-methylphenol	11.922	107	87777	9.352	ng	98
37) 2-Methylnaphthalene	12.280	142	211657	9.787	ng	98
38) 1-Methylnaphthalene	12.504	142	198014	9.783	ng	98
40) 1,2,4,5-Tetrachloroben...	12.651	216	102464	9.813	ng	99
41) Hexachlorocyclopentadiene	12.616	237	42768	7.742	ng	98
43) 2,4,6-Trichlorophenol	12.904	196	62645	9.360	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072823\
 Data File : BM041160.D
 Acq On : 28 Jul 2023 11:29
 Operator : MA/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/29/2023
 Supervised By :mohammad ahmed 07/29/2023

Quant Time: Jul 29 00:03:52 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 28 23:54:24 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.974	196	73138	9.442	ng	97
46) 1,1'-Biphenyl	13.298	154	257755	9.887	ng	99
47) 2-Chloronaphthalene	13.345	162	192927	9.911	ng	99
48) 2-Nitroaniline	13.569	65	53136	8.552	ng	99
49) Acenaphthylene	14.192	152	303860	9.671	ng	99
50) Dimethylphthalate	13.933	163	235376	9.730	ng	99
51) 2,6-Dinitrotoluene	14.063	165	46437	9.265	ng	95
52) Acenaphthene	14.533	154	185317	9.791	ng	97
53) 3-Nitroaniline	14.404	138	44307	9.807	ng	97
54) 2,4-Dinitrophenol	14.616	184	19018	11.044	ng	98
55) Dibenzofuran	14.868	168	306592	9.928	ng	100
56) 4-Nitrophenol	14.716	139	30771	7.366	ng	95
57) 2,4-Dinitrotoluene	14.863	165	54771	9.897	ng	# 87
58) Fluorene	15.521	166	234406	9.622	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.104	232	59088	9.420	ng	98
60) Diethylphthalate	15.304	149	223035	9.534	ng	99
61) 4-Chlorophenyl-phenyle...	15.521	204	115928	9.496	ng	99
62) 4-Nitroaniline	15.568	138	34529	10.049	ng	95
63) Azobenzene	15.815	77	241941	9.442	ng	99
65) 4,6-Dinitro-2-methylph...	15.621	198	28763	7.333	ng	89
66) n-Nitrosodiphenylamine	15.739	169	196123	9.739	ng	99
67) 4-Bromophenyl-phenylether	16.421	248	67812	9.552	ng	99
68) Hexachlorobenzene	16.527	284	77144	9.772	ng	99
69) Atrazine	16.698	200	53200	10.059	ng	94
70) Pentachlorophenol	16.880	266	48477	8.868	ng	98
71) Phenanthrene	17.268	178	356825	9.918	ng	100
72) Anthracene	17.362	178	338834	9.523	ng	100
73) Carbazole	17.645	167	308071	9.566	ng	99
74) Di-n-butylphthalate	18.203	149	321260	8.828	ng	99
75) Fluoranthene	19.298	202	385618	9.541	ng	99
77) Benzidine	19.503	184	77945	9.237	ng	98
78) Pyrene	19.656	202	405917	9.418	ng	100
80) Butylbenzylphthalate	20.568	149	133067	8.455	ng	98
81) Benzo(a)anthracene	21.415	228	400978	9.597	ng	99
82) 3,3'-Dichlorobenzidine	21.356	252	125559	9.210	ng	99
83) Chrysene	21.468	228	396218	9.805	ng	99
84) Bis(2-ethylhexyl)phtha...	21.339	149	188904	9.697	ng	99
85) Di-n-octyl phthalate	22.256	149	320400	8.055	ng	98
87) Indeno(1,2,3-cd)pyrene	26.250	276	487672	9.435	ng	100
88) Benzo(b)fluoranthene	23.080	252	399916	9.575	ng	99
89) Benzo(k)fluoranthene	23.127	252	400318	9.370	ng	99
90) Benzo(a)pyrene	23.697	252	372093	9.272	ng	99
91) Dibenzo(a,h)anthracene	26.262	278	405132	9.397	ng	98
92) Benzo(g,h,i)perylene	27.003	276	408012	9.672	ng	99

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

Manual Integrations

Quant Time: Jul 29 00:03:52 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 28 23:54:24 2023
Response via : Initial Calibration

