

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050824\
 Data File : BM045707.D
 Acq On : 08 May 2024 10:59
 Operator : MA/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 05/10/2024
 Supervised By :mohammad ahmed 05/10/2024

Quant Time: May 08 14:53:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM050824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 08 14:48:54 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.645	152	693017	20.000	ng	0.00
21) Naphthalene-d8	10.422	136	3212795	20.000	ng	0.00
39) Acenaphthene-d10	14.280	164	2042413	20.000	ng	0.00
64) Phenanthrene-d10	17.027	188	4190982	20.000	ng	0.00
76) Chrysene-d12	21.215	240	3835372	20.000	ng	0.00
86) Perylene-d12	23.415	264	4441008	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.251	112	804724	19.309	ng	0.00
7) Phenol-d6	6.816	99	1075125	19.025	ng	0.00
23) Nitrobenzene-d5	8.786	82	1064019	18.960	ng	0.00
42) 2,4,6-Tribromophenol	15.768	330	355652	17.776	ng	0.00
45) 2-Fluorobiphenyl	12.904	172	2575510	20.521	ng	0.00
79) Terphenyl-d14	19.674	244	4200348	18.720	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.228	88	157818	9.965	ng	99
3) Pyridine	3.604	79	446578	9.536	ng	98
4) n-Nitrosodimethylamine	3.522	42	235844	9.621	ng	# 97
6) Aniline	6.975	93	548575	9.792	ng	99
8) 2-Chlorophenol	7.210	128	451145	9.769	ng	97
9) Benzaldehyde	6.792	77	367150	9.854	ng	99
10) Phenol	6.839	94	553811	9.708	ng	99
11) bis(2-Chloroethyl)ether	7.081	93	474002	9.803	ng	98
12) 1,3-Dichlorobenzene	7.534	146	514085	9.916	ng	97
13) 1,4-Dichlorobenzene	7.681	146	523428	9.970	ng	97
14) 1,2-Dichlorobenzene	7.992	146	506079	10.078	ng	99
15) Benzyl Alcohol	7.881	79	381049	9.528	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.186	45	746828	10.017	ng	98
17) 2-Methylphenol	8.081	107	378817	9.774	ng	98
18) Hexachloroethane	8.716	117	198622	9.729	ng	96
19) n-Nitroso-di-n-propyla...	8.451	70	384828	9.920	ng	99
20) 3+4-Methylphenols	8.404	107	522754	9.658	ng	100
22) Acetophenone	8.457	105	751693	9.827	ng	98
24) Nitrobenzene	8.828	77	536823	9.610	ng	98
25) Isophorone	9.357	82	1094704	9.770	ng	99
26) 2-Nitrophenol	9.533	139	154407	10.690	ng	97
27) 2,4-Dimethylphenol	9.604	122	321137	9.478	ng	99
28) bis(2-Chloroethoxy)met...	9.845	93	667177	9.748	ng	99
29) 2,4-Dichlorophenol	10.063	162	389618	9.222	ng	97
30) 1,2,4-Trichlorobenzene	10.286	180	459556	9.617	ng	96
31) Naphthalene	10.469	128	1595437	9.827	ng	100
32) Benzoic acid	9.680	122	173022m	9.616	ng	
33) 4-Chloroaniline	10.575	127	622296	9.659	ng	99
34) Hexachlorobutadiene	10.769	225	262488	9.626	ng	97
35) Caprolactam	11.310	113	147280	9.234	ng	99
36) 4-Chloro-3-methylphenol	11.704	107	466394	9.385	ng	99
37) 2-Methylnaphthalene	12.086	142	1078163	9.708	ng	99
38) 1-Methylnaphthalene	12.310	142	1073493	9.778	ng	99
40) 1,2,4,5-Tetrachloroben...	12.457	216	474938	9.431	ng	99
41) Hexachlorocyclopentadiene	12.451	237	160659	9.032	ng	99
43) 2,4,6-Trichlorophenol	12.698	196	288582	8.864	ng	96

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM050824\
 Data File : BM045707.D
 Acq On : 08 May 2024 10:59
 Operator : MA/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 05/10/2024

Supervised By :mohammad ahmed 05/10/2024

Quant Time: May 08 14:53:48 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM050824.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed May 08 14:48:54 2024

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.763	196	324469	8.959	ng	97
46) 1,1'-Biphenyl	13.116	154	1381415	9.753	ng	100
47) 2-Chloronaphthalene	13.145	162	1083714	9.769	ng	99
48) 2-Nitroaniline	13.345	65	242801	10.346	ng	100
49) Acenaphthylene	13.998	152	1655575	9.772	ng	99
50) Dimethylphthalate	13.751	163	1350861	9.698	ng	100
51) 2,6-Dinitrotoluene	13.851	165	243521	8.723	ng	95
52) Acenaphthene	14.345	154	1046040m	9.788	ng	
53) 3-Nitroaniline	14.174	138	265774	8.495	ng	# 94
54) 2,4-Dinitrophenol	14.380	184	51881	10.003	ng	94
55) Dibenzofuran	14.680	168	1614549	9.951	ng	99
56) 4-Nitrophenol	14.480	139	209547	10.406	ng	97
57) 2,4-Dinitrotoluene	14.639	165	297576	10.311	ng	94
58) Fluorene	15.333	166	1347168	9.792	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.904	232	269715	9.030	ng	98
60) Diethylphthalate	15.127	149	1468769	9.724	ng	99
61) 4-Chlorophenyl-phenyle...	15.339	204	637672	9.781	ng	98
62) 4-Nitroaniline	15.333	138	302697	8.892	ng	96
63) Azobenzene	15.633	77	1469434	10.006	ng	99
65) 4,6-Dinitro-2-methylph...	15.404	198	80698	10.354	ng	96
66) n-Nitrosodiphenylamine	15.551	169	1119586	9.760	ng	99
67) 4-Bromophenyl-phenylether	16.233	248	384798	9.623	ng	99
68) Hexachlorobenzene	16.333	284	452456	9.705	ng	97
69) Atrazine	16.509	200	357215	10.269	ng	99
70) Pentachlorophenol	16.674	266	235777	8.077	ng	99
71) Phenanthrene	17.068	178	1966701	9.922	ng	100
72) Anthracene	17.156	178	1968871	9.953	ng	100
73) Carbazole	17.421	167	2006154	9.977	ng	100
74) Di-n-butylphthalate	18.027	149	2593656	10.446	ng	99
75) Fluoranthene	19.086	202	2366458	10.226	ng	99
77) Benzidine	19.274	184	616286	5.597	ng	99
78) Pyrene	19.445	202	2468907	10.003	ng	100
80) Butylbenzylphthalate	20.386	149	904682	10.357	ng	98
81) Benzo(a)anthracene	21.203	228	2431111	10.134	ng	98
82) 3,3'-Dichlorobenzidine	21.139	252	818839	9.159	ng	99
83) Chrysene	21.256	228	2229894	10.106	ng	99
84) Bis(2-ethylhexyl)phtha...	21.186	149	1618606	9.242	ng	99
85) Di-n-octyl phthalate	22.068	149	2741073	10.069	ng	99
87) Indeno(1,2,3-cd)pyrene	25.621	276	2773694	9.741	ng	100
88) Benzo(b)fluoranthene	22.762	252	2550842	9.806	ng	99
89) Benzo(k)fluoranthene	22.803	252	2336707	9.494	ng	99
90) Benzo(a)pyrene	23.321	252	2163439	9.470	ng	99
91) Dibenzo(a,h)anthracene	25.632	278	2298586	9.729	ng	99
92) Benzo(g,h,i)perylene	26.279	276	2272543	9.848	ng	99

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

