

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081224\
 Data File : BM047151.D
 Acq On : 12 Aug 2024 10:43
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 08/13/2024
 Supervised By :mohammad ahmed 08/14/2024

Quant Time: Aug 12 12:05:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:47:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.375	152	208548	20.000	ng	0.00
21) Naphthalene-d8	10.127	136	919233	20.000	ng	0.00
39) Acenaphthene-d10	14.033	164	654146	20.000	ng	0.00
64) Phenanthrene-d10	16.792	188	1480603	20.000	ng	0.00
76) Chrysene-d12	21.033	240	1414085	20.000	ng	0.00
86) Perylene-d12	23.727	264	1780647	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	892894	76.503	ng	0.00
7) Phenol-d6	6.581	99	1243186	77.303	ng	0.00
23) Nitrobenzene-d5	8.522	82	1421190	77.849	ng	0.00
42) 2,4,6-Tribromophenol	15.539	330	642308	84.682	ng	0.00
45) 2-Fluorobiphenyl	12.639	172	3128356	73.769	ng	0.00
79) Terphenyl-d14	19.462	244	5159238	78.178	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.034	88	186578	38.582	ng	99
3) Pyridine	3.410	79	570119m	40.204	ng	
4) n-Nitrosodimethylamine	3.328	42	247122	39.888	ng	96
6) Aniline	6.722	93	632369	41.602	ng	99
8) 2-Chlorophenol	6.951	128	507900	39.899	ng	99
9) Benzaldehyde	6.534	77	310310	34.973	ng	98
10) Phenol	6.610	94	611554	38.110	ng	97
11) bis(2-Chloroethyl)ether	6.822	93	552876	40.232	ng	97
12) 1,3-Dichlorobenzene	7.263	146	592678	39.181	ng	99
13) 1,4-Dichlorobenzene	7.410	146	606428	39.595	ng	100
14) 1,2-Dichlorobenzene	7.716	146	562769	38.944	ng	98
15) Benzyl Alcohol	7.622	79	475819	41.572	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.910	45	728000	41.284	ng	100
17) 2-Methylphenol	7.828	107	442039	41.290	ng	98
18) Hexachloroethane	8.428	117	245560	41.392	ng	97
19) n-Nitroso-di-n-propyla...	8.186	70	437413	40.379	ng	99
20) 3+4-Methylphenols	8.157	107	624560	41.730	ng	98
22) Acetophenone	8.192	105	809913	36.501	ng	98
24) Nitrobenzene	8.563	77	725805	39.483	ng	99
25) Isophorone	9.086	82	1256567	39.532	ng	99
26) 2-Nitrophenol	9.263	139	322749	39.645	ng	99
27) 2,4-Dimethylphenol	9.339	122	381773	40.034	ng	99
28) bis(2-Chloroethoxy)met...	9.575	93	808692	40.433	ng	99
29) 2,4-Dichlorophenol	9.792	162	563792	39.845	ng	98
30) 1,2,4-Trichlorobenzene	9.998	180	609035	38.035	ng	99
31) Naphthalene	10.180	128	1754672	38.318	ng	100
32) Benzoic acid	9.516	122	470853	42.499	ng	99
33) 4-Chloroaniline	10.304	127	737686	41.890	ng	99
34) Hexachlorobutadiene	10.463	225	356492	38.036	ng	99
35) Caprolactam	11.110	113	213722	43.494	ng	100
36) 4-Chloro-3-methylphenol	11.451	107	633755	42.056	ng	98
37) 2-Methylnaphthalene	11.804	142	1280920	39.286	ng	99
38) 1-Methylnaphthalene	12.027	142	1276021	39.599	ng	98
40) 1,2,4,5-Tetrachloroben...	12.186	216	652072	36.547	ng	99
41) Hexachlorocyclopentadiene	12.163	237	237406	38.105	ng	98
43) 2,4,6-Trichlorophenol	12.445	196	484531	38.207	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.516	196	543982	38.634	ng	96
46) 1,1'-Biphenyl	12.851	154	1726690	37.078	ng	99
47) 2-Chloronaphthalene	12.886	162	1348162	37.038	ng	100
48) 2-Nitroaniline	13.110	65	476203	41.761	ng	95
49) Acenaphthylene	13.745	152	2071676	38.612	ng	100
50) Dimethylphthalate	13.515	163	1822570	39.893	ng	100
51) 2,6-Dinitrotoluene	13.627	165	429849	40.249	ng	98
52) Acenaphthene	14.098	154	1315412	38.650	ng	100
53) 3-Nitroaniline	13.957	138	453216	42.350	ng	95
54) 2,4-Dinitrophenol	14.174	184	270595	42.275	ng	98
55) Dibenzofuran	14.439	168	2081182	38.714	ng	99
56) 4-Nitrophenol	14.292	139	403599	43.736	ng	91
57) 2,4-Dinitrotoluene	14.427	165	599300	42.042	ng	96
58) Fluorene	15.092	166	1770151	39.907	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.674	232	463368	40.042	ng	97
60) Diethylphthalate	14.892	149	1943451	41.657	ng	99
61) 4-Chlorophenyl-phenyle...	15.098	204	828620	39.239	ng	99
62) 4-Nitroaniline	15.133	138	456557	43.305	ng	96
63) Azobenzene	15.392	77	1844351	41.694	ng	99
65) 4,6-Dinitro-2-methylph...	15.198	198	363076	41.884	ng	98
66) n-Nitrosodiphenylamine	15.321	169	1538993	38.162	ng	99
67) 4-Bromophenyl-phenylether	15.992	248	544916	38.193	ng	95
68) Hexachlorobenzene	16.098	284	663319	38.819	ng	98
69) Atrazine	16.292	200	505062	41.619	ng	99
70) Pentachlorophenol	16.451	266	485836	40.561	ng	98
71) Phenanthrene	16.833	178	2906855	39.120	ng	100
72) Anthracene	16.927	178	2853566	39.468	ng	99
73) Carbazole	17.203	167	2959919	40.096	ng	100
74) Di-n-butylphthalate	17.792	149	3567201	41.529	ng	100
75) Fluoranthene	18.868	202	3649002	40.374	ng	100
77) Benzidine	19.074	184	1491270	62.186	ng	99
78) Pyrene	19.233	202	3867046	38.299	ng	99
80) Butylbenzylphthalate	20.174	149	1716787	41.664	ng	99
81) Benzo(a)anthracene	21.015	228	3703576	39.452	ng	100
82) 3,3'-Dichlorobenzidine	20.956	252	1244016	42.354	ng	99
83) Chrysene	21.074	228	3468062	38.938	ng	100
84) Bis(2-ethylhexyl)phtha...	20.974	149	2428011	41.549	ng	99
85) Di-n-octyl phthalate	22.009	149	4208330	42.369	ng	100
87) Indeno(1,2,3-cd)pyrene	26.726	276	4738171	41.167	ng	100
88) Benzo(b)fluoranthene	22.885	252	4156383	39.306	ng	100
89) Benzo(k)fluoranthene	22.944	252	3857875	38.624	ng	100
90) Benzo(a)pyrene	23.603	252	3650777	39.854	ng	98
91) Dibenzo(a,h)anthracene	26.779	278	3813537	40.942	ng	99
92) Benzo(g,h,i)perylene	27.662	276	4054202	41.922	ng	100

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

