

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011722\
 Data File : BM033757.D
 Acq On : 17 Jan 2022 19:49
 Operator : CG/JU
 Sample : PB142053BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :

BNA_M

ClientSampleId :

PB142053BS

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 01/18/2022
 Supervised By : Jagrut Upadhyay 01/18/2022

Quant Time: Jan 18 02:37:56 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM011722.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Jan 17 16:30:56 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.955	152	159262	20.000	ng	0.00
21) Naphthalene-d8	10.772	136	698430	20.000	ng	0.00
39) Acenaphthene-d10	14.595	164	466398	20.000	ng	0.00
64) Phenanthrene-d10	17.325	188	958998	20.000	ng	0.00
76) Chrysene-d12	21.495	240	903231	20.000	ng	0.00
86) Perylene-d12	23.906	264	889296	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.502	112	1421980	150.028	ng	0.00
7) Phenol-d6	7.119	99	1844946	138.168	ng	0.00
23) Nitrobenzene-d5	9.119	82	1226930	86.450	ng	0.00
42) 2,4,6-Tribromophenol	16.078	330	821800	129.965	ng	0.00
45) 2-Fluorobiphenyl	13.225	172	2848790	85.931	ng	0.00
79) Terphenyl-d14	19.942	244	4476142	82.181	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.337	88	131995	35.641	ng	# 92
3) Pyridine	3.749	79	354317	33.127	ng	92
4) n-Nitrosodimethylamine	3.655	42	228235	45.495	ng	# 72
6) Aniline	7.272	93	703751	45.403	ng	95
8) 2-Chlorophenol	7.519	128	566592	51.600	ng	87
9) Benzaldehyde	7.078	77	312182	45.730	ng	92
10) Phenol	7.143	94	703881	52.561	ng	93
11) bis(2-Chloroethyl)ether	7.372	93	499604	46.386	ng	93
12) 1,3-Dichlorobenzene	7.843	146	554631	45.664	ng	96
13) 1,4-Dichlorobenzene	7.990	146	572373	45.804	ng	97
14) 1,2-Dichlorobenzene	8.307	146	559884	45.946	ng	98
15) Benzyl Alcohol	8.196	79	547254	48.467	ng	92
16) 2,2'-oxybis(1-Chloropr...	8.490	45	703724	47.209	ng	96
17) 2-Methylphenol	8.402	107	487882	50.151	ng	95
18) Hexachloroethane	9.049	117	216248	46.801	ng	94
19) n-Nitroso-di-n-propyla...	8.772	70	424875	44.413	ng	# 88
20) 3+4-Methylphenols	8.731	107	664788	48.908	ng	94
22) Acetophenone	8.784	105	856188	46.430	ng	# 92
24) Nitrobenzene	9.160	77	634194	47.490	ng	92
25) Isophorone	9.696	82	1138842	45.563	ng	97
26) 2-Nitrophenol	9.878	139	305851	47.413	ng	# 84
27) 2,4-Dimethylphenol	9.943	122	550037	55.842	ng	94
28) bis(2-Chloroethoxy)met...	10.178	93	705868	44.323	ng	98
29) 2,4-Dichlorophenol	10.419	162	559638	49.692	ng	98
30) 1,2,4-Trichlorobenzene	10.631	180	550126	45.145	ng	98
31) Naphthalene	10.819	128	1728553	45.848	ng	99
32) Benzoic acid	10.096	122	390114	49.353	ng	89
33) 4-Chloroaniline	10.925	127	308313	19.652	ng	93
34) Hexachlorobutadiene	11.113	225	353459	44.263	ng	96
35) Caprolactam	11.719	113	181962m	45.371	ng	
36) 4-Chloro-3-methylphenol	12.054	107	665000	47.941	ng	99
37) 2-Methylnaphthalene	12.431	142	1276467	47.260	ng	98
38) 1-Methylnaphthalene	12.648	142	1182329	44.844	ng	98
40) 1,2,4,5-Tetrachloroben...	12.795	216	639919	47.415	ng	99
41) Hexachlorocyclopentadiene	12.784	237	937332	112.842	ng	98
43) 2,4,6-Trichlorophenol	13.031	196	454626	47.177	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.101	196	494204	47.623	ng	97
46) 1,1'-Biphenyl	13.431	154	1681907	47.649	ng	99
47) 2-Chloronaphthalene	13.472	162	1248736	46.362	ng	98
48) 2-Nitroaniline	13.672	65	425441	48.296	ng	# 83
49) Acenaphthylene	14.319	152	2049399	47.222	ng	99
50) Dimethylphthalate	14.054	163	1689059	45.964	ng	98
51) 2,6-Dinitrotoluene	14.166	165	363250	48.988	ng	86
52) Acenaphthene	14.660	154	1500759m	51.730	ng	
53) 3-Nitroaniline	14.489	138	214384	26.437	ng	90
54) 2,4-Dinitrophenol	14.695	184	408240	90.953	ng	88
55) Dibenzofuran	14.989	168	1944891	44.864	ng	97
56) 4-Nitrophenol	14.801	139	656908	98.719	ng	# 86
57) 2,4-Dinitrotoluene	14.948	165	514329	49.854	ng	85
58) Fluorene	15.636	166	1569158	45.924	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.213	232	430865	45.528	ng	99
60) Diethylphthalate	15.413	149	1764711	45.801	ng	100
61) 4-Chlorophenyl-phenyle...	15.631	204	799225	44.741	ng	97
62) 4-Nitroaniline	15.648	138	376961	44.262	ng	90
63) Azobenzene	15.919	77	1631102	46.043	ng	94
65) 4,6-Dinitro-2-methylph...	15.707	198	267761	42.629	ng	86
66) n-Nitrosodiphenylamine	15.842	169	1404950	47.073	ng	98
67) 4-Bromophenyl-phenylether	16.519	248	496072	45.528	ng	95
68) Hexachlorobenzene	16.642	284	556783	46.810	ng	96
69) Atrazine	16.789	200	536526	47.989	ng	99
70) Pentachlorophenol	16.983	266	753896	98.616	ng	99
71) Phenanthrene	17.372	178	2515147	47.112	ng	99
72) Anthracene	17.460	178	2574143	49.026	ng	98
73) Carbazole	17.724	167	2322007	45.312	ng	99
74) Di-n-butylphthalate	18.295	149	3061373	47.399	ng	100
75) Fluoranthene	19.377	202	2877086	48.129	ng	99
77) Benzidine	19.554	184	1390526	60.679	ng	99
78) Pyrene	19.736	202	2943322	44.108	ng	98
80) Butylbenzylphthalate	20.630	149	1370983	44.623	ng	95
81) Benzo(a)anthracene	21.477	228	2755686	46.096	ng	98
82) 3,3'-Dichlorobenzidine	21.407	252	667046	31.203	ng	99
83) Chrysene	21.530	228	2682388	46.975	ng	99
84) Bis(2-ethylhexyl)phtha...	21.407	149	2049711	46.170	ng	100
85) Di-n-octyl phthalate	22.336	149	3558180	49.879	ng	# 91
87) Indeno(1,2,3-cd)pyrene	26.436	276	2729145	46.160	ng	99
88) Benzo(b)fluoranthene	23.171	252	2807150	49.448	ng	99
89) Benzo(k)fluoranthene	23.218	252	2688260	47.283	ng	99
90) Benzo(a)pyrene	23.801	252	2633898	56.758	ng	99
91) Dibenzo(a,h)anthracene	26.447	278	2377085	47.595	ng	97
92) Benzo(g,h,i)perylene	27.206	276	2223883	46.793	ng	98

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

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