

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020922\
 Data File : BM034033.D
 Acq On : 09 Feb 2022 18:56
 Operator : CG/JU
 Sample : SP5814
 Misc : 8270/625 SPIKE
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP5814

Quant Time: Feb 10 04:06:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 09 17:52:36 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 02/10/2022
 Supervised By : Jagrut Upadhyay 02/11/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.900	152	140872	20.000	ng	0.00
21) Naphthalene-d8	10.715	136	535997	20.000	ng	# 0.00
39) Acenaphthene-d10	14.545	164	310757	20.000	ng	0.00
64) Phenanthrene-d10	17.270	188	558834	20.000	ng	# 0.00
76) Chrysene-d12	21.437	240	438871	20.000	ng	# 0.00
86) Perylene-d12	23.825	264	446426	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	0.000	112	0	0.000	ng	
7) Phenol-d6	0.000	99	0	0.000	ng	
23) Nitrobenzene-d5	0.000	82	0d	0.000	ng	
42) 2,4,6-Tribromophenol	0.000	330	0	0.000	ng	
45) 2-Fluorobiphenyl	0.000	172	0d	0.000	ng	
79) Terphenyl-d14	0.000	244	0d	0.000	ng	
Target Compounds						Qvalue
3) Pyridine	3.710	79	516911	49.053	ng	# 72
4) n-Nitrosodimethylamine	3.620	42	192673	50.690	ng	# 23
6) Aniline	7.224	93	570455	47.524	ng	# 89
8) 2-Chlorophenol	7.449	128	504937	58.320	ng	99
9) Benzaldehyde	7.021	77	424641	69.910	ng	98
10) Phenol	7.089	94	639416	58.474	ng	86
11) bis(2-Chloroethyl)ether	7.314	93	515211	57.742	ng	89
12) 1,3-Dichlorobenzene	7.787	146	603127	58.403	ng	96
13) 1,4-Dichlorobenzene	7.945	146	563570	56.531	ng	94
14) 1,2-Dichlorobenzene	8.260	146	556181	55.444	ng	96
15) Benzyl Alcohol	8.147	79	459101	56.192	ng	# 88
16) 2,2'-oxybis(1-Chloropr...	8.440	45	531381	55.266	ng	84
17) 2-Methylphenol	8.350	107	427478	59.918	ng	94
18) Hexachloroethane	8.981	117	220798	58.594	ng	90
19) n-Nitroso-di-n-propyla...	8.711	70	396096	57.879	ng	# 81
20) 3+4-Methylphenols	8.665	107	542028	55.084	ng	97
22) Acetophenone	8.733	105	736161	54.920	ng	# 82
24) Nitrobenzene	9.093	77	551157	54.070	ng	98
25) Isophorone	9.634	82	987849	59.296	ng	99
26) 2-Nitrophenol	9.814	139	255858	59.255	ng	91
27) 2,4-Dimethylphenol	9.882	122	487344	64.530	ng	95
28) bis(2-Chloroethoxy)met...	10.107	93	602343	54.702	ng	97
29) 2,4-Dichlorophenol	10.355	162	456116	57.627	ng	98
30) 1,2,4-Trichlorobenzene	10.580	180	540916	56.249	ng	98
31) Naphthalene	10.760	128	1581819	56.772	ng	99
32) Benzoic acid	10.017	122	258195	46.385	ng	93
33) 4-Chloroaniline	10.873	127	243780	22.340	ng	# 88
34) Hexachlorobutadiene	11.053	225	336306	54.942	ng	97
35) Caprolactam	11.639	113	116580	52.814	ng	# 76
36) 4-Chloro-3-methylphenol	11.999	107	486789	56.391	ng	90
37) 2-Methylnaphthalene	12.382	142	986164	55.483	ng	95
38) 1-Methylnaphthalene	12.585	142	994530	56.037	ng	98
40) 1,2,4,5-Tetrachloroben...	12.743	216	583979	56.371	ng	99
41) Hexachlorocyclopentadiene	12.720	237	850305	157.265	ng	96
43) 2,4,6-Trichlorophenol	12.968	196	333764	55.995	ng	100
44) 2,4,5-Trichlorophenol	13.058	196	360930	53.825	ng	# 84

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.373	154	1347962	56.653	ng	100
47) 2-Chloronaphthalene	13.418	162	1097705	57.477	ng	99
48) 2-Nitroaniline	13.621	65	287461	55.060	ng	# 67
49) Acenaphthylene	14.252	152	1540324	58.965	ng	99
50) Dimethylphthalate	14.004	163	1239390	55.658	ng	98
51) 2,6-Dinitrotoluene	14.117	165	277129	60.014	ng	# 74
52) Acenaphthene	14.612	154	1067242	59.576	ng	99
53) 3-Nitroaniline	14.432	138	132593	28.720	ng	97
54) 2,4-Dinitrophenol	14.657	184	271563	110.259	ng	# 59
55) Dibenzofuran	14.928	168	1350962	52.146	ng	91
56) 4-Nitrophenol	14.747	139	465793	116.107	ng	# 78
57) 2,4-Dinitrotoluene	14.905	165	349619	60.988	ng	# 57
58) Fluorene	15.581	166	1232854	54.076	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.175	232	286993	52.499	ng	# 34
60) Diethylphthalate	15.355	149	1284551	53.870	ng	98
61) 4-Chlorophenyl-phenyle...	15.581	204	676297	53.940	ng	95
62) 4-Nitroaniline	15.603	138	245324	51.472	ng	82
63) Azobenzene	15.874	77	1227332	54.810	ng	78
65) 4,6-Dinitro-2-methylph...	15.671	198	157479	49.214	ng	# 46
66) n-Nitrosodiphenylamine	15.783	169	966927	55.560	ng	99
67) 4-Bromophenyl-phenylether	16.482	248	321545	54.703	ng	# 77
68) Hexachlorobenzene	16.594	284	409280	55.717	ng	# 86
69) Atrazine	16.730	200	283553	53.918	ng	95
70) Pentachlorophenol	16.932	266	522382	115.639	ng	97
71) Phenanthrene	17.315	178	1777016	54.933	ng	99
72) Anthracene	17.405	178	1865811	58.941	ng	99
73) Carbazole	17.676	167	1619358	50.402	ng	100
74) Di-n-butylphthalate	18.239	149	2013234	58.151	ng	99
75) Fluoranthene	19.320	202	1643990	51.367	ng	94
77) Benzidine	19.500	184	953605	121.635	ng	99
78) Pyrene	19.680	202	1669728	59.655	ng	99
80) Butylbenzylphthalate	20.581	149	937389	50.166	ng	88
81) Benzo(a)anthracene	21.437	228	1603158	52.950	ng	99
82) 3,3'-Dichlorobenzidine	21.370	252	322910	39.718	ng	# 94
83) Chrysene	21.482	228	1771926m	53.833	ng	
84) Bis(2-ethylhexyl)phtha...	21.347	149	729388	49.268	ng	# 91
85) Di-n-octyl phthalate	22.271	149	1960692	55.131	ng	# 88
87) Indeno(1,2,3-cd)pyrene	26.325	276	1771301	54.554	ng	99
88) Benzo(b)fluoranthene	23.104	252	1797075m	59.140	ng	
89) Benzo(k)fluoranthene	23.149	252	1478394m	58.732	ng	
90) Benzo(a)pyrene	23.712	252	1576163	67.765	ng	# 97
91) Dibenzo(a,h)anthracene	26.325	278	1538919	57.682	ng	# 95
92) Benzo(g,h,i)perylene	27.091	276	1569320	58.795	ng	97

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

