

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020720\
 Data File : BG044342.D
 Acq On : 7 Feb 2020 18:17
 Operator : CG/JU
 Sample : L1408-02MSD
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SHORE-RD-PATH-BH-01-BMSD

Manual Integrations
 APPROVED

mohammad
 2/11/2020 8:26:08 AM

Quant Time: Feb 08 02:58:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG013020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 30 16:05:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	100025	20.00	ng	0.00
21) Naphthalene-d8	10.71	136	420044	20.00	ng	0.00
39) Acenaphthene-d10	14.55	164	266725	20.00	ng	0.00
64) Phenanthrene-d10	17.29	188	556346	20.00	ng	0.00
76) Chrysene-d12	21.54	240	471446	20.00	ng	0.00
87) Perylene-d12	24.63	264	540509	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	678102	119.64	ng	0.00
7) Phenol-d6	7.09	99	895973	117.72	ng	0.00
23) Nitrobenzene-d5	9.07	82	539363	72.49	ng	0.00
42) 2,4,6-Tribromophenol	16.04	330	335770	107.23	ng	0.00
45) 2-Fluorobiphenyl	13.17	172	1171325	73.54	ng	0.00
79) Terphenyl-d14	19.91	244	1575250	68.73	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.38	88	90033	35.356 ng	99
3) Pyridine	3.78	79	219209	32.311 ng	98
4) n-Nitrosodimethylamine	3.69	42	108787	38.373 ng	95
6) Aniline	7.24	93	199588	21.127 ng	97
8) 2-Chlorophenol	7.48	128	284467	45.669 ng	99
9) Benzaldehyde	7.05	77	126865	29.268 ng	98
10) Phenol	7.11	94	349261	46.369 ng	97
11) bis(2-Chloroethyl)ether	7.33	93	259745	40.336 ng	98
12) 1,3-Dichlorobenzene	7.81	146	291165	39.150 ng	96
13) 1,4-Dichlorobenzene	7.95	146	293711	39.874 ng	99
14) 1,2-Dichlorobenzene	8.27	146	275377	39.553 ng	99
15) Benzyl Alcohol	8.15	79	225661	41.880 ng	98
16) 2,2'-oxybis(1-Chloropropan	8.44	45	334507	38.380 ng	98
17) 2-Methylphenol	8.36	107	238459	44.372 ng	98
18) Hexachloroethane	9.00	117	105501	38.168 ng	98
19) n-Nitroso-di-n-propylamine	8.72	70	201780	39.781 ng	98
20) 3+4-Methylphenols	8.69	107	325758	43.984 ng	99
22) Acetophenone	8.73	105	379587	37.147 ng	99
24) Nitrobenzene	9.11	77	284960	39.232 ng	97
25) Isophorone	9.63	82	545824	39.360 ng	99
26) 2-Nitrophenol	9.83	139	159630	42.354 ng	99
27) 2,4-Dimethylphenol	9.89	122	259268	48.733 ng	100
28) bis(2-Chloroethoxy)methane	10.12	93	352993	38.157 ng	99
29) 2,4-Dichlorophenol	10.36	162	273554	42.523 ng	99
30) 1,2,4-Trichlorobenzene	10.58	180	278238	37.720 ng	98
31) Naphthalene	10.76	128	811885	39.839 ng	99
32) Benzoic acid	10.05	122	128724	40.059 ng	94
33) 4-Chloroaniline	10.87	127	78954	8.518 ng	98
34) Hexachlorobutadiene	11.05	225	166820	36.753 ng	99
35) Caprolactam	11.64	113	96585m	40.986 ng	
36) 4-Chloro-3-methylphenol	12.01	107	276536	41.000 ng	99
37) 2-Methylnaphthalene	12.37	142	595351	39.346 ng	99
38) 1-Methylnaphthalene	12.59	142	573651	40.213 ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.75	216	296335	37.141 ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.73	237	349037	91.171	ng	99
43) 2,4,6-Trichlorophenol	12.98	196	215021	41.694	ng	99
44) 2,4,5-Trichlorophenol	13.06	196	229736	43.614	ng	99
46) 1,1'-Biphenyl	13.38	154	753040	39.141	ng	99
47) 2-Chloronaphthalene	13.42	162	587602	38.381	ng	99
48) 2-Nitroaniline	13.62	65	168348	37.260	ng	99
49) Acenaphthylene	14.27	152	918894	40.922	ng	100
50) Dimethylphthalate	14.00	163	821064	42.824	ng	99
51) 2,6-Dinitrotoluene	14.12	165	167974	39.293	ng	98
52) Acenaphthene	14.62	154	563950	38.747	ng	97
53) 3-Nitroaniline	14.45	138	74884	15.833	ng	95
54) 2,4-Dinitrophenol	14.66	184	193812	76.845	ng	97
55) Dibenzofuran	14.95	168	873971	39.660	ng	100
56) 4-Nitrophenol	14.77	139	303120	87.495	ng	98
57) 2,4-Dinitrotoluene	14.91	165	229674	38.332	ng	99
58) Fluorene	15.60	166	683387	39.373	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.18	232	195893	41.172	ng	99
60) Diethylphthalate	15.37	149	723483	37.870	ng	99
61) 4-Chlorophenyl-phenylether	15.59	204	359986	38.252	ng	99
62) 4-Nitroaniline	15.61	138	159046	33.654	ng	98
63) Azobenzene	15.88	77	650777	38.867	ng	98
65) 4,6-Dinitro-2-methylphenol	15.68	198	143759	46.811	ng	98
66) n-Nitrosodiphenylamine	15.81	169	640327	41.070	ng	99
67) 4-Bromophenyl-phenylether	16.48	248	235008	38.772	ng	99
68) Hexachlorobenzene	16.61	284	257755	37.957	ng	98
69) Atrazine	16.75	200	239383	44.795	ng	97
70) Pentachlorophenol	16.95	266	288652	83.899	ng	97
71) Phenanthrene	17.34	178	1169769	43.421	ng	99
72) Anthracene	17.43	178	1136205	42.659	ng	99
73) Carbazole	17.69	167	980029	39.913	ng	100
74) Di-n-butylphthalate	18.26	149	1199394	39.528	ng	100
75) Fluoranthene	19.35	202	1529363	49.073	ng	97
77) Benzidine	19.52	184	258732	19.786	ng	97
78) Pyrene	19.71	202	1521971	50.269	ng	98
80) Butylbenzylphthalate	20.60	149	546085	39.579	ng	98
81) Benzo(a)anthracene	21.52	228	1369670	47.140	ng	99
82) 3,3'-Dichlorobenzidine	21.44	252	210331	18.652	ng	97
83) Chrysene	21.58	228	1267755	45.140	ng	99
84) Bis(2-ethylhexyl)phthalate	21.44	149	779996	41.072	ng	100
85) Di-n-octyl phthalate	22.61	149	1363039	41.482	ng	100
86) Indeno(1,2,3-cd)pyrene	28.14	276	1555119	42.442	ng	100
88) Benzo(b)fluoranthene	23.66	252	1459563	46.862	ng	100
89) Benzo(k)fluoranthene	23.72	252	1285816	42.358	ng	99
90) Benzo(a)pyrene	24.49	252	1316172	44.695	ng	99
91) Dibenzo(a,h)anthracene	28.20	278	1222338	41.941	ng	98
92) Benzo(g,h,i)perylene	29.23	276	1315455	45.088	ng	99

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

