

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072817\
 Data File : BM011053.D
 Acq On : 28 Jul 2017 21:56
 Operator : SJ/JU
 Sample : I4465-03MS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OK-03-072617-AMS

Manual Integrations
 APPROVED

Sohil
 7/31/2017 6:50:16 PM

Quant Time: Jul 31 17:05:17 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 26 17:29:06 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	158253	20.00	ng	0.00
21) Naphthalene-d8	10.15	136	684898	20.00	ng	0.00
38) Acenaphthene-d10	14.05	164	488062	20.00	ng	0.00
63) Phenanthrene-d10	16.80	188	1210723	20.00	ng	-0.01
75) Chrysene-d12	21.02	240	1162939	20.00	ng	0.00
86) Perylene-d12	23.13	264	137876	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.03	112	1416408	151.30	ng	0.00
7) Phenol-d6	6.59	99	1874585	140.93	ng	0.00
23) Nitrobenzene-d5	8.53	82	1701380	93.23	ng	0.00
41) 2,4,6-Tribromophenol	15.55	330	1126579	144.02	ng	0.00
44) 2-Fluorobiphenyl	12.66	172	3667325	94.24	ng	0.00
78) Terphenyl-d14	19.46	244	5226732	89.03	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.05	88	158988	37.825	ng	# 100
4) n-Nitrosodimethylamine	3.33	42	278060	47.525	ng	# 65
8) 2-Chlorophenol	6.96	128	414805	43.643	ng	# 78
9) Benzaldehyde	6.55	77	479309	55.694	ng	# 76
10) Phenol	6.62	94	622249	43.071	ng	# 93
11) bis(2-Chloroethyl)ether	6.84	93	572596	50.866	ng	# 95
12) 1,3-Dichlorobenzene	7.27	146	476352	41.124	ng	# 78
13) 1,4-Dichlorobenzene	7.42	146	480179	40.442	ng	# 88
14) 1,2-Dichlorobenzene	7.73	146	470888	41.480	ng	# 90
15) Benzyl Alcohol	7.63	79	222330	17.424	ng	# 83
16) 2,2'-oxybis(1-Chloropropan	7.93	45	460661	45.477	ng	# 74
17) 2-Methylphenol	7.84	107	421404	45.639	ng	# 72
18) Hexachloroethane	8.44	117	219987	42.486	ng	# 85
19) n-Nitroso-di-n-propylamine	8.20	70	540976	47.861	ng	# 94
20) 3+4-Methylphenols	8.17	107	568603	44.300	ng	# 80
22) Acetophenone	8.20	105	840070	40.918	ng	# 91
24) Nitrobenzene	8.57	77	850225	44.962	ng	# 86
25) Isophorone	9.10	82	1318283	43.327	ng	# 84
26) 2-Nitrophenol	9.27	139	247636	41.157	ng	# 14
27) 2,4-Dimethylphenol	9.35	122	452689	42.739	ng	# 62
28) bis(2-Chloroethoxy)methane	9.59	93	443264	26.121	ng	# 98
29) 2,4-Dichlorophenol	9.80	162	548646	46.096	ng	# 96
30) 1,2,4-Trichlorobenzene	10.02	180	633025	43.015	ng	# 98
31) Naphthalene	10.19	128	1459972	42.373	ng	# 99
32) Benzoic acid	9.50	122	320137	37.602	ng	# 68
33) 4-Chloroaniline	10.36	127	198865m	14.468	ng	# 99
34) Hexachlorobutadiene	10.48	225	491324	42.388	ng	# 99
35) Caprolactam	11.16	113	92324m	27.352	ng	# 76
36) 4-Chloro-3-methylphenol	11.45	107	632092	41.128	ng	# 85
37) 2-Methylnaphthalene	11.82	142	1212325	47.897	ng	# 100
39) 1,2,4,5-Tetrachlorobenzene	12.20	216	872682	41.618	ng	# 97
40) Hexachlorocyclopentadiene	12.18	237	1335847	109.332	ng	# 98
42) 2,4,6-Trichlorophenol	12.46	196	545115	42.951	ng	# 86
43) 2,4,5-Trichlorophenol	12.52	196	579885	44.362	ng	# 86

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1'-Biphenyl	12.86	154	1725320	40.883	nq	96
46) 2-Chloronaphthalene	12.90	162	1367611	43.983	nq	96
47) 2-Nitroaniline	13.12	65	325823	29.620	nq #	65
48) Acenaphthylene	13.76	152	1668743	35.961	nq	99
49) Dimethylphthalate	13.52	163	1729190	41.416	nq #	98
50) 2,6-Dinitrotoluene	13.63	165	361693	42.840	nq #	49
51) Acenaphthene	14.11	154	1245992	43.365	nq	99
52) 3-Nitroaniline	13.96	138	182942	24.970	nq #	36
53) 2,4-Dinitrophenol	14.17	184	513114	85.514	nq #	87
54) Dibenzofuran	14.45	168	2254201	48.420	nq #	91
55) 4-Nitrophenol	14.29	139	470945	73.162	nq	89
56) 2,4-Dinitrotoluene	14.43	165	503894	42.513	nq #	80
57) Fluorene	15.10	166	1826231	45.054	nq	97
58) 2,3,4,6-Tetrachlorophenol	14.68	232	615882	50.266	nq #	100
59) Diethylphthalate	14.90	149	1827652	42.863	nq	99
60) 4-Chlorophenyl-phenylether	15.11	204	1126045	46.012	nq	99
61) 4-Nitroaniline	15.14	138	216435	27.812	nq #	1
62) Azobenzene	15.40	77	2154754	47.217	nq	86
64) 4,6-Dinitro-2-methylphenol	15.19	198	336755	41.086	nq	94
65) n-Nitrosodiphenylamine	15.33	169	406777	11.740	nq	94
66) 4-Bromophenyl-phenylether	16.00	248	727458	46.739	nq #	89
67) Hexachlorobenzene	16.11	284	763974	43.438	nq #	87
69) Pentachlorophenol	16.46	266	940384	84.219	nq	96
70) Phenanthrene	16.84	178	2908586	45.880	nq	100
71) Anthracene	16.93	178	2839174	44.905	nq	98
72) Carbazole	17.21	167	332710	6.017	nq	98
73) Di-n-butylphthalate	17.80	149	2853486	42.378	nq #	97
74) Fluoranthene	18.87	202	3299624	42.000	nq	97
77) Pvrene	19.24	202	3019258	43.694	nq	99
79) Butylbenzylphthalate	20.18	149	1046066	42.572	nq #	69
80) Benzo(a)anthracene	21.00	228	2995430	45.110	nq	99
81) 3,3'-Dichlorobenzidine	20.96	252	81800	3.140	nq	96
82) Chrysene	21.06	228	2879004	45.168	nq	100
83) Bis(2-ethylhexyl)phthalate	20.97	149	1567739	43.370	nq #	96
84) Di-n-octyl phthalate	21.82	149	2573399	43.863	nq	99
85) Indeno(1,2,3-cd)pyrene	25.21	276	2452395	37.150	nq	96
87) Benzo(b)fluoranthene	22.51	252	3053393	345.052	nq #	92
88) Benzo(k)fluoranthene	22.55	252	2602750	306.876	nq #	96
89) Benzo(a)pyrene	23.04	252	2081468	259.949	nq #	95
90) Dibenzo(a,h)anthracene	25.22	278	2724870	367.086	nq #	91
91) Benzo(g,h,i)perylene	25.84	276	2024838	277.794	nq #	83

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

