

Data Path : Z:\HPCHEM1\BNA F\DATA\BF021318\
 Data File : BF102972.D
 Acq On : 13 Feb 2018 20:01
 Operator : SJ/JU
 Sample : J1537-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-01-COMP

Quant Time: Feb 14 00:10:40 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 00:38:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.03	152	442656	20.00	ng	0.15
21) Naphthalene-d8	8.50	136	325	20.00	ng	0.34
38) Acenaphthene-d10	10.37	164	2526	20.00	ng	0.45
63) Phenanthrene-d10	11.93	188	3968	20.00	ng	0.54
75) Chrysene-d12	14.67	240	1920	20.00	ng	0.64
86) Perylene-d12	16.24	264	8854	20.00	ng	0.74

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.50	112	1035334	35.52	ng	0.01
7) Phenol-d6	6.50	99	1242189	35.39	ng	0.00
23) Nitrobenzene-d5	0.00	82	0	0.00	ng	
41) 2,4,6-Tribromophenol	0.00	330	0	0.00	ng	
44) 2-Fluorobiphenyl	0.00	172	0	0.00	ng	
78) Terphenyl-d14	13.56	244	7455	91.94	ng	0.57

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
19) n-Nitroso-di-n-propylamine	7.44	70	99060	4.281	ng	# 81
22) Acetophenone	7.52	105	363	45.643	ng	# 8
24) Nitrobenzene	7.80	77	226	41.003	ng	# 38
25) Isophorone	8.00	82	158	14.646	ng	# 1
27) 2,4-Dimethylphenol	8.11	122	110	24.135	ng	# 67
28) bis(2-Chloroethoxy)methane	8.22	93	143	22.376	ng	# 23
31) Naphthalene	8.51	128	114	7.179	ng	# 1
32) Benzoic acid	8.11	122	110	37.781	ng	# 86
35) Caprolactam	8.97	113	637	442.705	ng	# 57
36) 4-Chloro-3-methylphenol	9.23	107	14887	3197.913	ng	# 1
37) 2-Methylnaphthalene	8.97	142	23332	2244.310	ng	# 95
45) 1,1'-Biphenyl	9.63	154	10270	48.271	ng	# 77
46) 2-Chloronaphthalene	9.62	162	526	3.140	ng	# 1
47) 2-Nitroaniline	9.89	65	845	18.638	ng	# 2
48) Acenaphthylene	10.22	152	3452	13.269	ng	# 1
50) 2,6-Dinitrotoluene	10.22	165	1872	49.958	ng	# 49
52) 3-Nitroaniline	10.32	138	124	2.689	ng	# 2
53) 2,4-Dinitrophenol	10.40	184	266	36.277	ng	# 54
54) Dibenzofuran	10.59	168	4754	21.683	ng	# 70
55) 4-Nitrophenol	10.70	139	2487	66.360	ng	# 87
56) 2,4-Dinitrotoluene	10.57	165	9471	179.459	ng	# 30
57) Fluorene	10.93	166	1216	7.623	ng	# 1
59) Diethylphthalate	10.77	149	613	3.152	ng	# 17
61) 4-Nitroaniline	10.99	138	248	5.651	ng	# 29
64) 4,6-Dinitro-2-methylphenol	11.00	198	600	39.281	ng	# 1
65) n-Nitrosodiphenylamine	11.06	169	3161	22.585	ng	# 33
68) Atrazine	11.73	200	722	17.486	ng	# 82
70) Phenanthrene	11.96	178	725	3.281	ng	# 1
71) Anthracene	12.04	178	758	3.372	ng	# 1
72) Carbazole	12.19	167	468	2.125	ng	# 1
73) Di-n-butylphthalate	12.54	149	4432	16.569	ng	# 1
74) Fluoranthene	13.20	202	1372	6.171	ng	# 1
76) Benzidine	13.32	184	855	9.689	ng	# 1
77) Pyrene	13.41	202	8881	58.987	ng	# 22

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79) Butylbenzylphthalate	14.07	149	2286	32.089	nq	91
81) 3,3'-Dichlorobenzidine	14.63	252	7557	168.791	nq #	1
82) Chrysene	14.72	228	879	7.221	nq #	1
83) Bis(2-ethylhexyl)phthalate	14.66	149	2877	31.194	nq #	74
84) Di-n-octyl phthalate	15.30	149	713	5.149	nq #	75
85) Indeno(1,2,3-cd)pyrene	17.47	276	189374	1875.849	nq	95
87) Benzo(b)fluoranthene	15.56	252	113918	198.451	nq #	86
88) Benzo(k)fluoranthene	15.56	252	113551	207.026	nq #	89

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

