

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE050917\
 Data File : BE092962.D
 Acq On : 9 May 2017 13:32
 Operator : SJ/MA
 Sample : PB98776BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB98776BS

Quant Time: May 09 17:13:41 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE050517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 09 15:34:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	1057	0.40	ng	0.00
7) Naphthalene-d8	10.53	136	3961	0.40	ng	0.00
13) Acenaphthene-d10	14.39	164	1717	0.40	ng	0.00
19) Phenanthrene-d10	17.13	188	3502	0.40	ng	0.00
26) Chrysene-d12	21.28	240	3839	0.40	ng	0.00
33) Perylene-d12	23.70	264	3436	0.40	ng	0.01

System Monitoring Compounds						
4) 2-Fluorophenol	5.36	112	815	0.35	ng	0.01
5) Phenol-d6	6.94	99	938	0.28	ng	0.00
8) Nitrobenzene-d5	8.90	82	840	0.31	ng	0.00
11) 2-Methylnaphthalene-d10	12.13	152	1711	0.29	ng	0.00
14) 2,4,6-Tribromophenol	15.88	330	92	0.34	ng	0.01
15) 2-Fluorobiphenyl	13.01	172	1976	0.25	ng	0.00
24) Fluoranthene-d10	19.15	212	2306	0.28	ng	0.00
28) Terphenyl-d14	19.74	244	1622	0.24	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	3.32	88	389	0.19	ng	# 78
3) n-Nitrosodimethylamine	3.61	42	457	0.29	ng	# 86
6) bis(2-Chloroethyl)ether	7.20	93	948	0.24	ng	99
9) Naphthalene	10.58	128	2703	0.26	ng	# 89
10) Hexachlorobutadiene	10.89	225	404	0.28	ng	93
12) 2-Methylnaphthalene	12.19	142	1604	0.25	ng	95
16) Acenaphthylene	14.10	152	7405	0.27	ng	99
17) Acenaphthene	14.45	154	1382	0.25	ng	96
18) Fluorene	15.44	166	1590	0.24	ng	100
20) Hexachlorobenzene	16.47	284	449	0.30	ng	# 100
21) Pentachlorophenol	16.81	266	192	0.66	ng	# 83
22) Phenanthrene	17.18	178	2690	0.27	ng	97
23) Anthracene	17.27	178	2083	0.28	ng	98
25) Fluoranthene	19.17	202	11742	0.28	ng	# 95
27) Pyrene	19.53	202	12322	0.25	ng	# 99
29) Benzo(a)anthracene	21.27	228	2256	0.25	ng	# 85
30) Chrysene	21.32	228	3144	0.27	ng	# 86
31) Bis(2-ethylhexyl)phthalate	21.20	149	737	0.30	ng	# 98
32) Indeno(1,2,3-cd)pyrene	26.24	276	13540	0.33	ng	99
34) Benzo(b)fluoranthene	22.96	252	3083	0.27	ng	# 88
35) Benzo(k)fluoranthene	23.01	252	3031	0.27	ng	# 83
36) Benzo(a)pyrene	23.58	252	2540	0.28	ng	# 79
37) Dibenzo(a,h)anthracene	26.27	278	2648	0.26	ng	# 83
38) Benzo(g,h,i)perylene	27.00	276	12716	0.29	ng	# 83

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

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