

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071615\
 Data File : BE090135.D
 Acq On : 17 Jul 2015 1:05
 Operator : TP/IZ
 Sample : PB84555BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB84555BL

Quant Time: Jul 17 13:15:58 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE071615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 17 12:59:41 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	11559	5.00	ng	0.00
6) Naphthalene-d8	10.37	136	51596	5.00	ng	0.00
12) Acenaphthene-d10	14.22	164	29074	5.00	ng	0.00
18) Phenanthrene-d10	16.96	188	65858	5.00	ng	0.00
25) Chrysene-d12	21.12	240	61806	5.00	ng	0.00
32) Perylene-d12	23.44	264	46412	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.24	112	15301	6.18	ng	0.00
5) Phenol-d6	6.79	99	22550	6.44	ng	0.00
7) Nitrobenzene-d5	8.75	82	13640	3.57	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	3111	3.86	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	30754	3.44	ng	0.00
27) Terphenyl-d14	19.58	244	41312	3.99	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
8) Nitrobenzene	8.79	77	869	0.22	ng	97
9) Naphthalene	10.41	128	2429	0.21	ng	94
10) Hexachlorobutadiene	10.71	225	277	0.14	ng	99
11) 2-Methylnaphthalene	12.04	142	1976	0.26	ng	99
15) Acenaphthylene	13.94	152	14451	0.28	ng	99
16) Acenaphthene	14.28	154	2199	0.29	ng	97
17) Fluorene	15.27	166	3085	0.32	ng	98
19) 4-Bromophenyl-phenylether	16.16	248	917	0.33	ng	93
20) Hexachlorobenzene	16.28	284	4040	0.33	ng	96
21) Pentachlorophenol	16.63	266	193	0.83	ng	92
22) Phenanthrene	17.00	178	5351	0.32	ng	99
23) Anthracene	17.10	178	4673	0.32	ng	100
24) Fluoranthene	19.01	202	5372	0.33	ng	100
26) Pyrene	19.37	202	5539	0.33	ng	100
28) Benzo(a)anthracene	21.11	228	3869	0.37	ng	98
29) Chrysene	21.16	228	6249	0.33	ng	98
30) Bis(2-ethylhexyl)phthalate	21.05	149	664	0.34	ng	95
31) Indeno(1,2,3-cd)pyrene	25.83	276	1278	0.40	ng	98
33) Benzo(b)fluoranthene	22.73	252	1390	0.41	ng	# 89
34) Benzo(k)fluoranthene	22.78	252	2863	0.37	ng	# 92
35) Benzo(a)pyrene	23.33	252	1597	0.56	ng	# 86
36) Dibenzo(a,h)anthracene	25.86	278	811	0.41	ng	# 81
37) Benzo(g,h,i)perylene	26.57	276	1772	0.54	ng	# 91

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

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