

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070815\
 Data File : BE090071.D
 Acq On : 8 Jul 2015 22:44
 Operator : TP/IZ
 Sample : G2811-06MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X1SS0720006-150625MS

Quant Time: Jul 09 01:21:06 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 30 07:01:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	20855	5.00	ng	-0.01
6) Naphthalene-d8	10.37	136	94061	5.00	ng	-0.01
12) Acenaphthene-d10	14.22	164	51310	5.00	ng	-0.01
18) Phenanthrene-d10	16.96	188	112948	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	123339	5.00	ng	-0.01
32) Perylene-d12	23.44	264	115156	5.00	ng	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Fluorophenol	5.24	112	29387	6.13	ng	0.00
5) Phenol-d6	6.79	99	48735	7.27	ng	0.00
7) Nitrobenzene-d5	8.75	82	29376	3.93	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	13145	8.74	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	61280	3.97	ng	-0.01
27) Terphenyl-d14	19.58	244	87374	4.25	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.24	88	6974	2.44	ng	98
3) n-Nitrosodimethylamine	3.53	42	11201	2.89	ng	91
8) Nitrobenzene	8.79	77	27906	3.52	ng	96
9) Naphthalene	10.41	128	72066	3.39	ng	99
10) Hexachlorobutadiene	10.72	225	11151	3.34	ng	100
11) 2-Methylnaphthalene	12.03	142	49487	3.48	ng	99
15) Acenaphthylene	13.94	152	330659	3.57	ng	99
16) Acenaphthene	14.29	154	48776	3.66	ng	96
17) Fluorene	15.27	166	64480	3.55	ng	99
19) 4-Bromophenyl-phenylether	16.16	248	17670	3.68	ng	94
20) Hexachlorobenzene	16.28	284	72980	3.65	ng	97
21) Pentachlorophenol	16.63	266	20083	7.85	ng	95
22) Phenanthrene	17.00	178	116331	4.02	ng	99
23) Anthracene	17.10	178	104783	3.96	ng	99
24) Fluoranthene	19.00	202	172524	5.34	ng	99
26) Pyrene	19.37	202	175514	5.97	ng	98
28) Benzo(a)anthracene	21.10	228	155673	5.05	ng	99
29) Chrysene	21.16	228	154234	4.82	ng	99
30) Bis(2-ethylhexyl)phthalate	21.05	149	74951	5.27	ng	97
31) Indeno(1,2,3-cd)pyrene	25.84	276	146520	4.78	ng	99
33) Benzo(b)fluoranthene	22.73	252	162904	6.41	ng	99
34) Benzo(k)fluoranthene	22.78	252	137415	4.84	ng	99
35) Benzo(a)pyrene	23.33	252	144671	6.17	ng	100
36) Dibenzo(a,h)anthracene	25.86	278	102419	4.25	ng	99
37) Benzo(g,h,i)perylene	26.57	276	129428	5.30	ng	98

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

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