

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071615\
 Data File : BE090149.D
 Acq On : 17 Jul 2015 10:30
 Operator : TP/IZ
 Sample : G3015-01MSD 10X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 STOCKPILE-7-15-15MSD

Quant Time: Jul 17 13:58:55 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	11394	5.00	ng	0.00
6) Naphthalene-d8	10.37	136	53119	5.00	ng	0.00
12) Acenaphthene-d10	14.22	164	30377	5.00	ng	0.00
18) Phenanthrene-d10	16.96	188	66942	5.00	ng	0.00
25) Chrysene-d12	21.13	240	70072	5.00	ng	0.00
32) Perylene-d12	23.45	264	69858	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	2101	0.86	ng	0.00
5) Phenol-d6	6.80	99	3230	0.94	ng	0.00
7) Nitrobenzene-d5	8.75	82	1876	0.48	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	863	1.29	ng	0.00
14) 2-Fluorobiphenyl	12.84	172	4266	0.46	ng	0.00
27) Terphenyl-d14	19.58	244	7273	0.62	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.24	88	572	0.37	ng	# 85
3) n-Nitrosodimethylamine	3.53	42	661	0.37	ng	# 78
8) Nitrobenzene	8.79	77	1747	0.42	ng	99
9) Naphthalene	10.41	128	67251	5.59	ng	99
10) Hexachlorobutadiene	10.71	225	825	0.41	ng	99
11) 2-Methylnaphthalene	12.03	142	19640	2.55	ng	100
15) Acenaphthylene	13.94	152	161654	2.99	ng	99
16) Acenaphthene	14.28	154	96953	12.41	ng	94
17) Fluorene	15.27	166	108994	10.68	ng	98
19) 4-Bromophenyl-phenylether	16.16	248	1233	0.43	ng	# 63
20) Hexachlorobenzene	16.28	284	5237	0.42	ng	95
21) Pentachlorophenol	16.63	266	986	1.61	ng	94
22) Phenanthrene	17.01	178	1587758	94.22	ng	97
23) Anthracene	17.10	178	505835m	33.87	ng	
24) Fluoranthene	19.02	202	2410051	144.74	ng	95
26) Pyrene	19.38	202	2228298	117.60	ng	# 88
28) Benzo(a)anthracene	21.11	228	1275713	107.40	ng	93
29) Chrysene	21.16	228	1179117	67.32	ng	94
30) Bis(2-ethylhexyl)phthalate	21.05	149	5527	2.21	ng	# 88
31) Indeno(1,2,3-cd)pyrene	25.86	276	570103	122.07	ng	95
33) Benzo(b)fluoranthene	22.75	252	1388850	273.47	ng	98
34) Benzo(k)fluoranthene	22.79	252	560797m	25.86	ng	
35) Benzo(a)pyrene	23.35	252	1001349	67.36	ng	97
36) Dibenzo(a,h)anthracene	25.87	278	164136m	11.34	ng	
37) Benzo(g,h,i)perylene	26.61	276	584642	38.94	ng	94

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

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