

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072315\
 Data File : BM002020.D
 Acq On : 23 Jul 2015 17:22
 Operator : TP/UM
 Sample : G3000-15RE
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C02C2RE

Quant Time: Jul 24 01:59:42 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM071315.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 24 01:29:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	77078	20.00	ng/ul	0.00
18) Naphthalene-d8	10.40	136	290750	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.28	164	160574	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.03	188	342789	20.00	ng/ul	0.00
77) Chrysene-d12	21.23	240	387470	20.00	ng/ul	0.00
85) Perylene-d12	23.44	264	393955	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.09	96	9063	5.75	ng/uL	0.00
5) Phenol-d5	6.79	99	202245	32.13	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.95	67	112474	32.80	ng/ul	0.00
9) 2-Chlorophenol-d4	7.15	132	178134	35.65	ng/ul	0.00
13) 4-Methylphenol-d8	8.33	113	164875	31.90	ng/ul	0.00
19) Nitrobenzene-d5	8.77	128	78275	36.04	ng/ul	0.00
22) 2-Nitrophenol-d4	9.49	143	87537	36.38	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.03	165	163854	33.68	ng/ul	0.00
29) 4-Chloroaniline-d4	10.55	131	156192	31.20	ng/ul	0.00
43) Dimethylphthalate-d6	13.69	166	416093	32.13	ng/ul	0.00
46) Acenaphthylene-d8	13.97	160	435774	27.98	ng/ul	0.00
51) 4-Nitrophenol-d4	14.49	143	66660	30.39	ng/ul	0.00
57) Fluorene-d10	15.27	176	269476	24.35	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.41	200	21258	9.58	ng/ul	0.00
70) Anthracene-d10	17.13	188	393774	24.47	ng/ul	0.00
78) Pyrene-d10	19.43	212	372767	21.06	ng/ul	0.00
89) Benzo(a)pyrene-d12	23.30	264	360929	18.12	ng/ul	0.00

Target Compounds

						Qvalue
44) Dimethylphthalate	13.74	163	219235	16.64	ng/ul	99
75) Di-n-butylphthalate	18.00	149	90539	4.54	ng/ul	97
76) Fluoranthene	19.09	202	36448	1.68	ng/ul#	97
79) Pyrene	19.46	202	38733	1.67	ng/ul	97
84) Chrysene	21.26	228	27270	1.26	ng/ul#	85
87) Benzo(b)fluoranthene	22.78	252	55635	2.30	ng/ul#	62
90) Benzo(a)pyrene	23.35	252	29511	1.30	ng/ul#	54
91) Indeno(1,2,3-cd)pyrene	25.67	276	30165	1.27	ng/ul#	93
93) Benzo(g,h,i)perylene	26.35	276	30178	1.54	ng/ul#	86

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

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