

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061116\
 Data File : BM005737.D
 Acq On : 12 Jun 2016 01:02
 Operator : UM/SJ
 Sample : H3411-25MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XR2MSD

Manual Integrations
 APPROVED

umangi
 6/13/2016 7:40:47 PM

Quant Time: Jun 12 02:22:59 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jun 11 23:08:56 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	101466	20.00	ng/ul	0.00
7) Naphthalene-d8	10.53	136	532136	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.39	164	336593	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.13	188	742629	20.00	ng/ul	0.00
29) Chrysene-d12	21.32	240	628614	20.00	ng/ul	0.00
34) Perylene-d12	23.57	264	526348	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.19	96	1766	0.75	ng/uL	0.00
3) Phenol-d5	6.93	99	73063	8.34	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.08	67	54624	10.38	ng/ul	0.00
5) 2-Chlorophenol-d4	7.28	132	68296	9.65	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	34076	4.61	ng/ul	0.00
8) Nitrobenzene-d5	8.90	128	34729	9.06	ng/ul	0.00
9) 2-Nitrophenol-d4	9.62	143	41835	9.65	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	75442	9.59	ng/ul	0.00
12) 4-Chloroaniline-d4	10.68	131	56967	6.00	ng/ul	0.00
16) Dimethylphthalate-d6	13.80	166	286738	10.33	ng/ul	0.00
17) Acenaphthylene-d8	14.08	160	374068	11.01	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	24232	5.97	ng/ul	0.00
21) Fluorene-d10	15.38	176	261074	11.11	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.52	200	22438	5.84	ng/ul	0.00
26) Anthracene-d10	17.23	188	394914	11.29	ng/ul	0.00
30) Pyrene-d10	19.53	212	439829	14.56	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.43	264	285087	11.88	ng/ul	0.00

Target Compounds

					Qvalue
19) Acenaphthene	14.45	153	271128	11.98 ng/ul	98
25) Phenanthrene	17.17	178	123111	3.03 ng/ul	100
28) Fluoranthene	19.19	202	297860	6.24 ng/ul	98
31) Pyrene	19.56	202	865782	22.63 ng/ul	100
32) Benzo(a)anthracene	21.30	228	130368	3.59 ng/ul	99
33) Chrysene	21.35	228	131370	3.67 ng/ul	97
35) Benzo(b)fluoranthene	22.89	252	154025	4.93 ng/ul	98
36) Benzo(k)fluoranthene	22.93	252	48368m	1.53 ng/ul	
38) Benzo(a)pyrene	23.47	252	102148	3.44 ng/ul	99
39) Indeno(1,2,3-cd)pyrene	25.84	276	72287	2.49 ng/ul	95
41) Benzo(g,h,i)perylene	26.54	276	74700	3.08 ng/ul	97

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

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