

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060616\
 Data File : BM005665.D
 Acq On : 06 Jun 2016 12:11
 Operator : UM/SJ
 Sample : H3412-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XS6

Manual Integrations
 APPROVED

sohil
 6/7/2016 7:13:29 PM

Quant Time: Jun 07 04:33:27 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 07 04:25:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.76	152	86015	20.00	ng/ul	0.00
7) Naphthalene-d8	10.55	136	400749	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	212692	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	393132	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	335940	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	353711	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.22	96	2709	1.32	ng/uL	0.00
3) Phenol-d5	6.94	99	87016	11.51	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.09	67	71292	15.49	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	82794	13.27	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	32548	5.44	ng/ul	0.00
8) Nitrobenzene-d5	8.92	128	42824	13.92	ng/ul	0.00
9) 2-Nitrophenol-d4	9.64	143	42710	12.97	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.18	165	69666	11.90	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	85319	12.27	ng/ul	0.00
16) Dimethylphthalate-d6	13.81	166	247539	15.14	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	318715	14.99	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	11417	4.88	ng/ul	0.00
21) Fluorene-d10	15.39	176	214705	15.29	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	8518	4.60	ng/ul	0.00
26) Anthracene-d10	17.24	188	279378	14.96	ng/ul	0.00
30) Pyrene-d10	19.54	212	269249	15.95	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	226937	13.94	ng/ul	0.00

Target Compounds

						Qvalue
25) Phenanthrene	17.19	178	27432	1.26	ng/ul	98
28) Fluoranthene	19.20	202	66768	3.01	ng/ul	98
31) Pyrene	19.56	202	59501	2.81	ng/ul	98
32) Benzo(a)anthracene	21.31	228	43862	2.23	ng/ul	97
33) Chrysene	21.36	228	48613	2.61	ng/ul	98
35) Benzo(b)fluoranthene	22.91	252	72237	3.39	ng/ul	97
36) Benzo(k)fluoranthene	22.94	252	20164m	1.01	ng/ul	
38) Benzo(a)pyrene	23.49	252	37356	1.82	ng/ul#	96
39) Indeno(1,2,3-cd)pyrene	25.87	276	29643	1.22	ng/ul#	93
41) Benzo(g,h,i)perylene	26.57	276	27802	1.36	ng/ul#	96

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

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