

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060616\
 Data File : BM005680.D
 Acq On : 06 Jun 2016 21:28
 Operator : UM/SJ
 Sample : H3412-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XS8

Manual Integrations
 APPROVED

sohil
 6/7/2016 7:13:57 PM

Quant Time: Jun 07 05:04:23 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	86379	20.00	ng/ul	0.00
7) Naphthalene-d8	10.55	136	359892	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	191929	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	403442	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	411548	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	429751	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.22	96	2840	1.38	ng/uL	0.00
3) Phenol-d5	6.94	99	84325	11.10	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.09	67	71481	15.47	ng/ul	0.00
5) 2-Chlorophenol-d4	7.30	132	83946	13.40	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	23907	3.98	ng/ul	0.00
8) Nitrobenzene-d5	8.92	128	41778	15.12	ng/ul	0.00
9) 2-Nitrophenol-d4	9.64	143	43181	14.60	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.18	165	64332	12.24	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	51964	8.32	ng/ul	0.00
16) Dimethylphthalate-d6	13.81	166	219726	14.89	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	285086	14.86	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	11037m	5.22	ng/ul	0.01
21) Fluorene-d10	15.39	176	193323	15.25	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	10126	5.33	ng/ul	0.00
26) Anthracene-d10	17.24	188	268658	14.02	ng/ul	0.00
30) Pyrene-d10	19.54	212	293159	14.17	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	279973	14.16	ng/ul	0.00

Target Compounds

						Qvalue
18) Acenaphthylene	14.12	152	35139	1.77	ng/ul	96
25) Phenanthrene	17.19	178	45445	2.04	ng/ul	97
27) Anthracene	17.28	178	27419	1.21	ng/ul	99
28) Fluoranthene	19.20	202	143067	6.29	ng/ul	98
31) Pyrene	19.57	202	162721	6.27	ng/ul	97
32) Benzo(a)anthracene	21.31	228	112000	4.64	ng/ul	97
33) Chrysene	21.36	228	133364	5.85	ng/ul	96
35) Benzo(b)fluoranthene	22.91	252	273699	10.56	ng/ul	98
36) Benzo(k)fluoranthene	22.94	252	75376m	3.10	ng/ul	
38) Benzo(a)pyrene	23.49	252	153774	6.16	ng/ul	96
39) Indeno(1,2,3-cd)pyrene	25.87	276	129638m	4.39	ng/ul	
40) Dibenzo(a,h)anthracene	25.87	278	38853m	1.57	ng/ul	
41) Benzo(g,h,i)perylene	26.58	276	114763	4.63	ng/ul#	94

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

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