

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081016\
 Data File : BM007023.D
 Acq On : 11 Aug 2016 05:00
 Operator : UM/SJ
 Sample : H4302-06DL 4X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleID :
 DA0H0DL

Manual Integrations
 APPROVED

umangi
 8/11/2016 7:55:00 AM

Quant Time: Aug 11 06:54:18 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080916.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 10 01:30:54 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	388616	20.00	ng/ul	0.00
7) Naphthalene-d8	10.60	136	1712708	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.45	164	1062791	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.19	188	2532370	20.00	ng/ul	0.00
29) Chrysene-d12	21.37	240	2497964	20.00	ng/ul	0.00
34) Perylene-d12	23.65	264	2114333	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.33	96	2474	0.29	ng/uL	0.00
3) Phenol-d5	6.97	99	116993	3.38	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.15	67	72362	3.74	ng/ul	0.00
5) 2-Chlorophenol-d4	7.35	132	96309	3.67	ng/ul	0.00
6) 4-Methylphenol-d8	8.51	113	84168	2.87	ng/ul	0.00
8) Nitrobenzene-d5	8.97	128	45344	3.47	ng/ul	0.00
9) 2-Nitrophenol-d4	9.69	143	47793	3.19	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.22	165	108740	3.72	ng/ul	0.00
12) 4-Chloroaniline-d4	10.74	131	82219	2.40	ng/ul	0.00
16) Dimethylphthalate-d6	13.86	166	372623	4.10	ng/ul	-0.01
17) Acenaphthylene-d8	14.14	160	452886	4.26	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	36275	2.22	ng/ul	-0.01
21) Fluorene-d10	15.44	176	339825	4.40	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.55	200	30620	2.23	ng/ul	0.00
26) Anthracene-d10	17.29	188	536059	4.54	ng/ul	0.00
30) Pyrene-d10	19.58	212	619867	5.46	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.51	264	446196	4.59	ng/ul	0.00

Target Compounds

						Qvalue
11) Naphthalene	10.65	128	169964	1.98	ng/ul	97
19) Acenaphthene	14.51	153	221472	3.12	ng/ul	96
22) Fluorene	15.49	166	132560	1.57	ng/ul	99
25) Phenanthrene	17.23	178	3209477	23.66	ng/ul	99
27) Anthracene	17.32	178	613004	4.37	ng/ul	100
28) Fluoranthene	19.25	202	5948700	34.81	ng/ul	98
31) Pyrene	19.62	202	4926970	34.00	ng/ul	98
32) Benzo(a)anthracene	21.36	228	2749193	18.69	ng/ul	99
33) Chrysene	21.41	228	2531116	18.96	ng/ul	98
35) Benzo(b)fluoranthene	22.97	252	3559095	27.10	ng/ul	98
36) Benzo(k)fluoranthene	23.01	252	1212276m	9.77	ng/ul	
38) Benzo(a)pyrene	23.55	252	2309378	19.11	ng/ul	97
39) Indeno(1,2,3-cd)pyrene	25.96	276	1947389	16.25	ng/ul	97
40) Dibenzo(a,h)anthracene	25.96	278	470908m	4.75	ng/ul	
41) Benzo(g,h,i)perylene	26.66	276	1907899	19.37	ng/ul	98

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

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