

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM073116\
 Data File : BM006810.D
 Acq On : 01 Aug 2016 04:23
 Operator : UM/SJ
 Sample : H4190-16
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA0D3

Manual Integrations
 APPROVED

sohil
 8/1/2016 7:08:30 PM

Quant Time: Aug 01 08:11:44 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Jul 30 01:37:26 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	225289	20.00	ng/ul	0.00
7) Naphthalene-d8	10.39	136	951384	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.26	164	509217	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.01	188	1103956	20.00	ng/ul	0.00
29) Chrysene-d12	21.22	240	1140062	20.00	ng/ul	0.00
34) Perylene-d12	23.43	264	1206507	20.00	ng/ul	0.01

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.13	96	8617	1.55	ng/uL	0.00
3) Phenol-d5	6.79	99	297249	15.51	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	6.95	67	204581	17.23	ng/ul	0.00
5) 2-Chlorophenol-d4	7.14	132	255461	16.68	ng/ul	0.00
6) 4-Methylphenol-d8	8.32	113	208756	14.04	ng/ul	0.00
8) Nitrobenzene-d5	8.76	128	125684	17.45	ng/ul	0.00
9) 2-Nitrophenol-d4	9.48	143	138535	18.09	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.02	165	238202	16.46	ng/ul	0.00
12) 4-Chloroaniline-d4	10.53	131	279942	17.56	ng/ul	0.00
16) Dimethylphthalate-d6	13.67	166	713931	17.88	ng/ul	0.00
17) Acenaphthylene-d8	13.95	160	926943	18.19	ng/ul	0.00
20) 4-Nitrophenol-d4	14.50	143	93271	13.45	ng/ul	0.01
21) Fluorene-d10	15.26	176	631236	17.68	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.40	200	61204	10.68	ng/ul	0.00
26) Anthracene-d10	17.11	188	944143	18.01	ng/ul	0.00
30) Pyrene-d10	19.42	212	1006960	19.38	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.29	264	990823	17.85	ng/ul	0.00

Target Compounds

						Qvalue
25) Phenanthrene	17.06	178	136501	2.20	ng/ul	98
28) Fluoranthene	19.08	202	366007	5.02	ng/ul	98
31) Pyrene	19.44	202	317469	4.60	ng/ul	98
32) Benzo(a)anthracene	21.20	228	192880	2.79	ng/ul	93
33) Chrysene	21.25	228	238486	3.74	ng/ul	96
35) Benzo(b)fluoranthene	22.76	252	324592	4.49	ng/ul#	96
36) Benzo(k)fluoranthene	22.80	252	100403m	1.44	ng/ul	
38) Benzo(a)pyrene	23.33	252	216258	3.07	ng/ul	95
39) Indeno(1,2,3-cd)pyrene	25.63	276	157132	1.93	ng/ul	95
41) Benzo(g,h,i)perylene	26.30	276	169115	2.42	ng/ul#	93

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

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