

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG051916\
 Data File : BG021965.D
 Acq On : 19 May 2016 16:09
 Operator : UM/SJ
 Sample : H3092-21
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9XD3

Manual Integrations
 APPROVED

sohil
 5/20/2016 7:16:26 PM

Quant Time: May 20 06:59:28 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG051816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 20 05:44:47 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	37933	20.00	ng/ul	0.00
7) Naphthalene-d8	10.98	136	198177	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.78	164	110028	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.53	188	207694	20.00	ng/ul	0.00
29) Chrysene-d12	21.81	240	178939	20.00	ng/ul	0.00
34) Perylene-d12	25.15	264	172180	20.00	ng/ul	0.01

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.54	96	2424	2.93	ng/uL	0.00
3) Phenol-d5	7.32	99	39376	11.76	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.48	67	25887	14.51	ng/ul	0.00
5) 2-Chlorophenol-d4	7.69	132	38011	13.82	ng/ul	0.00
6) 4-Methylphenol-d8	8.87	113	15652	5.25	ng/ul	0.00
8) Nitrobenzene-d5	9.33	128	20221	14.54	ng/ul	0.00
9) 2-Nitrophenol-d4	10.05	143	22644	15.26	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.60	165	32619	12.44	ng/ul	0.00
12) 4-Chloroaniline-d4	11.12	131	36187	17.52	ng/ul	0.00
16) Dimethylphthalate-d6	14.18	166	128299	15.12	ng/ul	0.00
17) Acenaphthylene-d8	14.48	160	176419	16.14	ng/ul	0.00
20) 4-Nitrophenol-d4	15.00	143	9728m	6.52	ng/ul	0.02
21) Fluorene-d10	15.78	176	114315	14.97	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.89	200	9420	9.14	ng/ul	0.00
26) Anthracene-d10	17.63	188	151803m	15.86	ng/ul	0.00
30) Pyrene-d10	19.91	212	150025	15.89	ng/ul	0.00
37) Benzo(a)pyrene-d12	24.91	264	123513	16.15	ng/ul	0.00

Target Compounds

						Qvalue
25) Phenanthrene	17.57	178	23332	2.10	ng/ul#	94
28) Fluoranthene	19.57	202	59386	5.15	ng/ul#	94
31) Pyrene	19.94	202	51200	4.15	ng/ul	96
32) Benzo(a)anthracene	21.79	228	28546	2.88	ng/ul	96
33) Chrysene	21.86	228	31014m	3.21	ng/ul	
35) Benzo(b)fluoranthene	24.09	252	46750	4.72	ng/ul#	90
36) Benzo(k)fluoranthene	24.14	252	13729m	1.41	ng/ul	
38) Benzo(a)pyrene	24.98	252	28950	3.06	ng/ul#	89
39) Indeno(1,2,3-cd)pyrene	28.95	276	21166m	1.97	ng/ul	
41) Benzo(g,h,i)perylene	30.15	276	24061m	2.64	ng/ul	

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

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