

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110215\
 Data File : BG019444.D
 Acq On : 2 Nov 2015 22:36
 Operator : UM/NP
 Sample : G4202-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BCY38

Manual Integrations
 APPROVED

Mmdadoda
 11/3/2015 5:57:43 PM

Quant Time: Nov 03 01:08:32 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG102815.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 03 00:57:47 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.18	152	24385	20.00	ng/ul	0.00
18) Naphthalene-d8	11.02	136	94881	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.82	164	80420	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.56	188	249282	20.00	ng/ul	0.00
78) Chrysene-d12	21.84	240	356257	20.00	ng/ul	0.00
86) Perylene-d12	25.21	264	333294	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.50	96	668	1.31	ng/uL	0.00
5) Phenol-d5	7.34	99	15742	7.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.50	67	54645	38.51	ng/ul	0.00
9) 2-Chlorophenol-d4	7.71	132	41152	29.45	ng/ul	0.00
13) 4-Methylphenol-d8	8.89	113	31456	17.64	ng/ul	0.00
19) Nitrobenzene-d5	9.36	128	29111	41.75	ng/ul	0.00
22) 2-Nitrophenol-d4	10.09	143	34744	41.97	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.63	165	65577	36.52	ng/ul	0.00
29) 4-Chloroaniline-d4	11.15	131	5856	3.22	ng/ul	0.00
44) Dimethylphthalate-d6	14.22	166	302218	45.58	ng/ul	0.00
47) Acenaphthylene-d8	14.52	160	294829	40.15	ng/ul	0.00
52) 4-Nitrophenol-d4	15.01	143	10038	9.53	ng/ul	0.00
58) Fluorene-d10	15.80	176	244828	39.91	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.91	200	48086	29.80	ng/ul	0.00
71) Anthracene-d10	17.66	188	463987	40.83	ng/ul	0.00
79) Pyrene-d10	19.93	212	608869	39.05	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.97	264	681684	40.76	ng/ul	0.00

Target Compounds

					Ovalue
81) Butylbenzylphthalate	20.83	149	7892	1.21	ng/ul 95
82) 3,3'-Dichlorobenzidine	21.73	252	8383	1.20	ng/ul 97
84) Bis(2-ethylhexyl)phthalate	21.71	149	23441	2.54	ng/ul# 89
87) Di-n-octyl phthalate	22.97	149	35660	2.36	ng/ul 100
88) Benzo(b)fluoranthene	24.13	252	28715	1.46	ng/ul# 96
89) Benzo(k)fluoranthene	24.20	252	26324	1.39	ng/ul# 90
91) Benzo(a)pyrene	25.05	252	26706	1.41	ng/ul# 91
92) Indeno(1,2,3-cd)pyrene	29.04	276	33259m	1.56	ng/ul
93) Dibenzo(a,h)anthracene	29.11	278	27137m	1.53	ng/ul
94) Benzo(g,h,i)perylene	30.26	276	29970m	1.86	ng/ul

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

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