

Data Path : Z:\HPCHEM1\BNA G\DATA\BG013117\
 Data File : BG025734.D
 Acq On : 1 Feb 2017 2:23
 Operator : SJ/MA
 Sample : PB96531BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK31

Manual Integrations
 APPROVED

mohammad
 2/1/2017 4:56:51 PM

Quant Time: Feb 01 04:41:38 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG011817.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 01 04:33:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	81916	20.00	ng/ul	0.00
18) Naphthalene-d8	11.00	136	343075	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.79	164	264698	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.54	188	582498	20.00	ng/ul	0.00
75) Chrysene-d12	21.84	240	816421	20.00	ng/ul	0.00
83) Perylene-d12	25.22	264	809929	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.50	96	10484m	6.40	ng/uL	0.00
5) Phenol-d5	7.33	99	249672	35.56	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.48	67	185447	40.90	ng/ul	0.00
9) 2-Chlorophenol-d4	7.70	132	181110	36.57	ng/ul	0.00
13) 4-Methylphenol-d8	8.88	113	211340	35.41	ng/ul	0.00
19) Nitrobenzene-d5	9.35	128	96652	39.74	ng/ul	0.00
22) 2-Nitrophenol-d4	10.08	143	115990	39.97	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.63	165	183206m	32.00	ng/ul	0.00
29) 4-Chloroaniline-d4	11.15	131	265505	41.49	ng/ul	0.00
43) Dimethylphthalate-d6	14.19	166	739064	39.54	ng/ul	0.00
46) Acenaphthylene-d8	14.49	160	831734	41.19	ng/ul	0.00
51) 4-Nitrophenol-d4	15.04	143	95753m	37.38	ng/ul	0.00
57) Fluorene-d10	15.79	176	639086	36.27	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.94	200	119148	33.16	ng/ul	0.00
70) Anthracene-d10	17.64	188	1043692	42.08	ng/ul	0.00
76) Pyrene-d10	19.92	212	1373480	40.43	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.98	264	1410647	41.44	ng/ul	0.00

Target Compounds	Ovalue

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

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