

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG080319\
 Data File : BG041920.D
 Acq On : 3 Aug 2019 22:23
 Operator : HP/JU
 Sample : PB121669BL
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK69

Quant Time: Aug 04 02:57:19 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Aug 04 02:31:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	68058	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	287754	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.69	164	192786	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.44	188	419210	20.00	ng/ul	0.00
77) Chrysene-d12	21.73	240	408645	20.00	ng/ul	0.00
85) Perylene-d12	25.00	264	431284	20.00	ng/ul	-0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.42	96	9465	6.08	ng/uL	0.00
5) Phenol-d5	7.19	99	202007	37.82	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.35	67	112788	37.76	ng/ul	0.00
9) 2-Chlorophenol-d4	7.56	132	178970	39.19	ng/ul	0.00
13) 4-Methylphenol-d8	8.74	113	170766	38.08	ng/ul	0.00
19) Nitrobenzene-d5	9.20	128	89242	41.20	ng/ul	0.00
22) 2-Nitrophenol-d4	9.93	143	107180	42.62	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.48	165	201134	39.61	ng/ul	0.00
29) 4-Chloroaniline-d4	10.99	131	246938	43.83	ng/ul	0.00
43) Dimethylphthalate-d6	14.09	166	590025	39.46	ng/ul	0.00
46) Acenaphthylene-d8	14.39	160	689589	40.37	ng/ul	0.00
51) 4-Nitrophenol-d4	14.88	143	90948	36.23	ng/ul	0.00
57) Fluorene-d10	15.69	176	530886	39.91	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.80	200	67193	24.98	ng/ul	0.00
70) Anthracene-d10	17.54	188	826020	41.07	ng/ul	0.00
78) Pyrene-d10	19.83	212	909641	39.89	ng/ul	-0.01
89) Benzo(a)pyrene-d12	24.78	264	971078	43.10	ng/ul	0.00

Target Compounds	Ovalue

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

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