

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN022220\
 Data File : BN009814.D
 Acq On : 21 Feb 2020 19:07
 Operator : CG/JU
 Sample : PB126856BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK56

Quant Time: Feb 21 22:03:15 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 19 04:46:48 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	126526	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.14	136	544918	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.03	164	353029	20.00	ng/ul	-0.02
62) Phenanthrene-d10	16.79	188	750511	20.00	ng/ul	-0.02
78) Chrysene-d12	21.01	240	710732	20.00	ng/ul	-0.01
86) Perylene-d12	23.10	264	717921	20.00	ng/ul	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.04	96	16499	5.36	ng/uL	-0.01
5) Phenol-d5	6.60	99	284129	26.40	ng/ul	-0.02
7) Bis-(2-Chloroethyl)ether-d	6.75	67	212974	28.68	ng/ul	-0.02
9) 2-Chlorophenol-d4	6.93	132	227565	28.09	ng/ul	-0.02
13) 4-Methylphenol-d8	8.10	113	230838	26.96	ng/ul	-0.02
19) Nitrobenzene-d5	8.53	128	114603	28.75	ng/ul	-0.02
22) 2-Nitrophenol-d4	9.25	143	127436	29.22	ng/ul	-0.02
26) 2,4-Dichlorophenol-d3	9.78	165	211684	25.25	ng/ul	-0.02
29) 4-Chloroaniline-d4	10.29	131	311570	33.08	ng/ul	-0.02
44) Dimethylphthalate-d6	13.46	166	767963	29.12	ng/ul	-0.02
47) Acenaphthylene-d8	13.72	160	880723	28.36	ng/ul	-0.02
52) 4-Nitrophenol-d4	14.27	143	102221	21.23	ng/ul	-0.02
58) Fluorene-d10	15.04	176	663815	29.16	ng/ul	-0.02
63) 4,6-Dinitro-2-methylphenol	15.19	200	104883	22.50	ng/ul	-0.02
71) Anthracene-d10	16.89	188	1057705	30.74	ng/ul	-0.02
79) Pyrene-d10	19.20	212	1168725	31.46	ng/ul	-0.02
90) Benzo(a)pyrene-d12	22.97	264	1151432	32.12	ng/ul	-0.03

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

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