

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121919\
 Data File : BM024177.D
 Acq On : 19 Dec 2019 16:59
 Operator : JU
 Sample : PB125361BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK61

Quant Time: Dec 20 04:05:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 17 15:54:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	275219	20.00	ng/ul	-0.02
18) Naphthalene-d8	10.20	136	1198135	20.00	ng/ul	-0.02
36) Acenaphthene-d10	14.10	164	755754	20.00	ng/ul	-0.01
62) Phenanthrene-d10	16.86	188	1585737	20.00	ng/ul	-0.01
78) Chrysene-d12	21.07	240	1490195	20.00	ng/ul	0.00
86) Perylene-d12	23.20	264	1809111	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.99	96	46840	7.52	ng/uL	-0.01
5) Phenol-d5	6.63	99	781920	30.01	ng/ul	-0.02
7) Bis-(2-Chloroethyl)ether-d	6.79	67	556601	35.61	ng/ul	-0.01
9) 2-Chlorophenol-d4	6.98	132	636919	34.27	ng/ul	-0.02
13) 4-Methylphenol-d8	8.16	113	649446	30.49	ng/ul	-0.01
19) Nitrobenzene-d5	8.59	128	315890	36.69	ng/ul	-0.01
22) 2-Nitrophenol-d4	9.31	143	356729	37.69	ng/ul	-0.01
26) 2,4-Dichlorophenol-d3	9.85	165	598383	32.52	ng/ul	-0.01
29) 4-Chloroaniline-d4	10.36	131	898534	40.20	ng/ul	-0.02
44) Dimethylphthalate-d6	13.53	166	2129988	38.11	ng/ul	-0.01
47) Acenaphthylene-d8	13.79	160	2482580	37.07	ng/ul	-0.01
52) 4-Nitrophenol-d4	14.35	143	258137	25.73	ng/ul	-0.01
58) Fluorene-d10	15.10	176	1812135	36.99	ng/ul	-0.01
63) 4,6-Dinitro-2-methylphenol	15.26	200	287602	29.73	ng/ul	0.00
71) Anthracene-d10	16.96	188	2695475	37.36	ng/ul	-0.01
79) Pyrene-d10	19.27	212	2899391	40.85	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.07	264	3404874	37.89	ng/ul	0.00

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

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