

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM121219\
 Data File : BM024081.D
 Acq On : 13 Dec 2019 08:33
 Operator : JU
 Sample : PB125245BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK45

Quant Time: Dec 13 13:39:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 07 04:26:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	510649	20.00	ng/ul	0.00
18) Naphthalene-d8	10.25	136	2184374	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.13	164	1386600	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.89	188	2987636	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	2689778	20.00	ng/ul	0.00
86) Perylene-d12	23.24	264	2561987	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.03	96	78608	6.85	ng/uL	0.00
5) Phenol-d5	6.67	99	1473636	32.93	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.83	67	949461	36.48	ng/ul	0.00
9) 2-Chlorophenol-d4	7.01	132	1170009	33.82	ng/ul	0.00
13) 4-Methylphenol-d8	8.19	113	1192645	33.77	ng/ul	0.00
19) Nitrobenzene-d5	8.63	128	589225	34.75	ng/ul	0.00
22) 2-Nitrophenol-d4	9.34	143	647189	33.59	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.88	165	1139644	31.23	ng/ul	0.00
29) 4-Chloroaniline-d4	10.39	131	1603214	40.03	ng/ul	0.00
44) Dimethylphthalate-d6	13.56	166	3744238	34.75	ng/ul	0.00
47) Acenaphthylene-d8	13.82	160	4432188	35.10	ng/ul	0.00
52) 4-Nitrophenol-d4	14.36	143	559028	30.86	ng/ul	0.00
58) Fluorene-d10	15.13	176	3187785	35.16	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.28	200	505959	26.98	ng/ul	0.00
71) Anthracene-d10	16.99	188	4949406	34.69	ng/ul	0.00
79) Pyrene-d10	19.29	212	5340662	38.43	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.11	264	4820924	35.60	ng/ul	0.00

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

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