

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122319\
 Data File : BM024217.D
 Acq On : 23 Dec 2019 15:59
 Operator : JU
 Sample : PB125443BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK43

Quant Time: Dec 23 17:09:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM121719MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Dec 21 00:34:28 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.43	152	214941	20.00	ng/ul	0.00
18) Naphthalene-d8	10.20	136	939738	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.10	164	597487	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.86	188	1279312	20.00	ng/ul	0.00
78) Chrysene-d12	21.07	240	1193166	20.00	ng/ul	0.00
86) Perylene-d12	23.20	264	1410837	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.00	96	37433	7.69	ng/uL	0.00
5) Phenol-d5	6.63	99	653104	32.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.79	67	473789	38.82	ng/ul	0.00
9) 2-Chlorophenol-d4	6.97	132	521153	35.91	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	532789	32.03	ng/ul	0.00
19) Nitrobenzene-d5	8.59	128	256648	38.01	ng/ul	0.00
22) 2-Nitrophenol-d4	9.30	143	277720	37.41	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.85	165	484559	33.57	ng/ul	0.00
29) 4-Chloroaniline-d4	10.36	131	712992	40.67	ng/ul	0.00
44) Dimethylphthalate-d6	13.53	166	1786350	40.43	ng/ul	0.00
47) Acenaphthylene-d8	13.79	160	2030208	38.35	ng/ul	0.00
52) 4-Nitrophenol-d4	14.35	143	202340	25.51	ng/ul	0.00
58) Fluorene-d10	15.10	176	1496784	38.64	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.26	200	208433	26.71	ng/ul	0.00
71) Anthracene-d10	16.96	188	2312195	39.72	ng/ul	0.00
79) Pyrene-d10	19.26	212	2489072	43.79	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.07	264	2814110	40.16	ng/ul	0.01

Target Compounds	Ovalue

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

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