

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM010920\
 Data File : BM024422.D
 Acq On : 09 Jan 2020 13:42
 Operator : JU
 Sample : PB125829BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK29

Quant Time: Jan 09 14:17:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM010820MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 09 01:07:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	270609	20.00	ng/ul	0.00
18) Naphthalene-d8	10.26	136	1082998	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.15	164	633061	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.89	188	1326903	20.00	ng/ul	0.00
78) Chrysene-d12	21.10	240	1263366	20.00	ng/ul	0.00
86) Perylene-d12	23.23	264	1482776	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.10	96	50973	8.13	ng/uL	0.00
5) Phenol-d5	6.69	99	790091	37.39	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.85	67	506209	40.85	ng/ul	0.00
9) 2-Chlorophenol-d4	7.04	132	690392	38.65	ng/ul	0.00
13) 4-Methylphenol-d8	8.20	113	637320	37.55	ng/ul	0.00
19) Nitrobenzene-d5	8.63	128	337242	41.52	ng/ul	0.00
22) 2-Nitrophenol-d4	9.35	143	312570	40.15	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.89	165	602868	35.29	ng/ul	0.00
29) 4-Chloroaniline-d4	10.39	131	903367	46.89	ng/ul	0.00
44) Dimethylphthalate-d6	13.57	166	1959885	43.10	ng/ul	0.00
47) Acenaphthylene-d8	13.83	160	2394756	40.79	ng/ul	0.00
52) 4-Nitrophenol-d4	14.35	143	321946	40.15	ng/ul	0.00
58) Fluorene-d10	15.14	176	1666202	41.34	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.26	200	157473	26.49	ng/ul	0.00
71) Anthracene-d10	16.99	188	2569459	41.35	ng/ul	0.00
79) Pyrene-d10	19.29	212	2872496	42.56	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.10	264	3179300	41.99	ng/ul	0.00

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

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