

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030921\
 Data File : BN013883.D
 Acq On : 09 Mar 2021 09:41
 Operator : CG/JU
 Sample : PB134929BL
 Misc : SOM-MDL-W-BL-04
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK29

Quant Time: Mar 09 10:14:23 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN030421MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 09 09:50:00 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	195093	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	810521	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.35	164	474541	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.09	188	977195	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	956216	20.00	ng/ul	0.00
86) Perylene-d12	23.50	264	980520	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.22	96	31249	6.36	ng/uL	0.00
5) Phenol-d5	6.88	99	553529	35.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	349628	35.90	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	441932	36.63	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	427530	35.78	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	204799	37.76	ng/ul	0.00
22) 2-Nitrophenol-d4	9.59	143	204304	39.43	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.12	165	396737	35.05	ng/ul	0.00
29) 4-Chloroaniline-d4	10.63	131	599674	42.79	ng/ul	0.00
44) Dimethylphthalate-d6	13.76	166	1329678	41.73	ng/ul	0.00
47) Acenaphthylene-d8	14.05	160	1616753	38.37	ng/ul	0.00
52) 4-Nitrophenol-d4	14.55	143	210574	35.23	ng/ul	0.00
58) Fluorene-d10	15.35	176	1120106	39.81	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.46	200	141405	31.13	ng/ul	0.00
71) Anthracene-d10	17.19	188	1785493	41.81	ng/ul	0.00
79) Pyrene-d10	19.48	212	1961505	42.17	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.36	264	2058928	42.18	ng/ul	0.00

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

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