

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030921\
 Data File : BN013888.D
 Acq On : 09 Mar 2021 12:43
 Operator : CG/JU
 Sample : PB134930BL
 Misc : SOM-MDL-S-BL-05
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK30

Quant Time: Mar 09 13:19:59 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	158129	20.00	ng/ul	0.00
18) Naphthalene-d8	10.50	136	650277	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.35	164	382998	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.09	188	783511	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	769597	20.00	ng/ul	0.00
86) Perylene-d12	23.49	264	793553	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	25346	6.36	ng/uL	0.00
5) Phenol-d5	6.88	99	439424	34.36	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	283775	35.95	ng/ul	0.00
9) 2-Chlorophenol-d4	7.25	132	355538	36.35	ng/ul	0.00
13) 4-Methylphenol-d8	8.42	113	338837	34.99	ng/ul	0.00
19) Nitrobenzene-d5	8.87	128	159752	36.71	ng/ul	0.00
22) 2-Nitrophenol-d4	9.58	143	155547	37.42	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.12	165	316245	34.83	ng/ul	0.00
29) 4-Chloroaniline-d4	10.63	131	484636	43.10	ng/ul	0.00
44) Dimethylphthalate-d6	13.76	166	1069846	41.60	ng/ul	0.00
47) Acenaphthylene-d8	14.05	160	1314868	38.66	ng/ul	0.00
52) 4-Nitrophenol-d4	14.54	143	160998	33.38	ng/ul	0.00
58) Fluorene-d10	15.35	176	920020	40.51	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.46	200	108219	29.72	ng/ul	0.00
71) Anthracene-d10	17.19	188	1466130	42.82	ng/ul	0.00
79) Pyrene-d10	19.49	212	1568346	41.90	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.35	264	1683543	42.61	ng/ul	0.00

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

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