

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP011121\
 Data File : BP004494.D
 Acq On : 11 Jan 2021 14:28
 Operator : CG/JU
 Sample : PB134017BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK17

Quant Time: Jan 11 15:09:06 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SOM-EPA-BP121820MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 24 16:46:25 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	226230	20.00	ng/ul	0.00
18) Naphthalene-d8	10.61	136	847076	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.47	164	515136	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.23	188	1057754	20.00	ng/ul	0.00
78) Chrysene-d12	21.32	240	1023086	20.00	ng/ul	0.00
86) Perylene-d12	23.68	264	1034768	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.30	96	40703	7.48	ng/uL	-0.01
5) Phenol-d5	6.99	99	676646	34.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.15	67	356674	32.38	ng/ul	0.00
9) 2-Chlorophenol-d4	7.35	132	583661	36.33	ng/ul	0.00
13) 4-Methylphenol-d8	8.52	113	531264	33.96	ng/ul	0.00
19) Nitrobenzene-d5	8.98	128	267828	36.53	ng/ul	0.00
22) 2-Nitrophenol-d4	9.70	143	306919	38.39	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.23	165	543822	36.68	ng/ul	0.00
29) 4-Chloroaniline-d4	10.75	131	715614	42.57	ng/ul	0.00
44) Dimethylphthalate-d6	13.87	166	1549246	37.57	ng/ul	0.00
47) Acenaphthylene-d8	14.16	160	2021250	39.47	ng/ul	0.00
52) 4-Nitrophenol-d4	14.67	143	224735	30.17	ng/ul	0.00
58) Fluorene-d10	15.46	176	1335294	37.02	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.60	200	194627	28.52	ng/ul	0.00
71) Anthracene-d10	17.33	188	2068457	39.58	ng/ul	0.00
79) Pyrene-d10	19.57	212	2272295	41.57	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.53	264	2410073	42.28	ng/ul	0.00

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

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