

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP021820\
 Data File : BP001764.D
 Acq On : 18 Feb 2020 17:48
 Operator : CG/JU
 Sample : PB126811BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK11

Quant Time: Feb 19 10:57:22 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP021820MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 19 02:37:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	350511	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	1400827	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.29	164	860839	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.05	188	1896667	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1996324	20.00	ng/ul	0.00
86) Perylene-d12	23.35	264	2173243	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.18	96	54539	6.46	ng/uL	0.00
5) Phenol-d5	6.83	99	849124	32.06	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	508735	33.06	ng/ul	0.00
9) 2-Chlorophenol-d4	7.19	132	737808	34.47	ng/ul	0.00
13) 4-Methylphenol-d8	8.36	113	662566	31.54	ng/ul	0.00
19) Nitrobenzene-d5	8.80	128	327117	34.50	ng/ul	0.00
22) 2-Nitrophenol-d4	9.52	143	318191	37.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.06	165	667183	32.72	ng/ul	0.00
29) 4-Chloroaniline-d4	10.56	131	1003251	41.59	ng/ul	0.00
44) Dimethylphthalate-d6	13.71	166	2225865	36.85	ng/ul	0.00
47) Acenaphthylene-d8	13.98	160	2496556	33.56	ng/ul	0.00
52) 4-Nitrophenol-d4	14.49	143	299598	28.48	ng/ul	0.00
58) Fluorene-d10	15.29	176	1919524	35.33	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.40	200	183795	24.58	ng/ul	0.00
71) Anthracene-d10	17.14	188	2954694	35.36	ng/ul	0.00
79) Pyrene-d10	19.39	212	3418586	33.37	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.21	264	3677135	35.03	ng/ul	0.00

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

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