

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
 Data File : BP015669.D
 Acq On : 23 Jun 2023 13:49
 Operator : MA/JU
 Sample : PB153415BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK415

Quant Time: Jun 23 22:44:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 22 22:01:49 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.081	152	116280	20.000	ng/u1	0.00
20) Naphthalene-d8	10.910	136	461334	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.734	164	257592	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.492	188	473994	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	418839	20.000	ng/u1	0.00
88) Perylene-d12	24.151	264	525088	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	20868	6.646	ng/uL	0.00
4) Pyridine-d5	3.840	84	271535	33.290	ng/u1	0.00
7) Phenol-d5	7.246	99	356235	33.254	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.405	67	232296	36.309	ng/u1	0.00
11) 2-Chlorophenol-d4	7.611	132	295228	38.012	ng/u1	0.00
15) 4-Methylphenol-d8	8.805	113	294709	33.521	ng/u1	0.00
21) Nitrobenzene-d5	9.263	128	142473	39.336	ng/u1	0.00
24) 2-Nitrophenol-d4	9.993	143	165794	40.083	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	262610	34.805	ng/u1	0.00
31) 4-Chloroaniline-d4	11.057	131	362120	33.436	ng/u1	0.00
46) Dimethylphthalate-d6	14.140	166	807114	39.734	ng/u1	0.00
49) Acenaphthylene-d8	14.428	160	907032	40.837	ng/u1	0.00
54) 4-Nitrophenol-d4	14.969	143	84402	23.203	ng/u1	0.02
60) Fluorene-d10	15.728	176	629751	36.765	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	87782	29.234	ng/u1	0.01
73) Anthracene-d10	17.598	188	874919	40.569	ng/u1	0.00
81) Pyrene-d10	19.822	212	957863	42.729	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.980	264	1090165	41.378	ng/u1	0.01

Target Compounds Qvalue

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

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