

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015851.D
 Acq On : 30 Jun 2023 19:05
 Operator : MA/JU
 Sample : PB153574BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK574

Quant Time: Jun 30 22:35:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 30 22:30:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.054	152	195707	20.000	ng/ul	0.00
20) Naphthalene-d8	10.884	136	878010	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.707	164	521302	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.466	188	1061838	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.565	240	741664	20.000	ng/ul	# 0.00
88) Perylene-d12	24.100	264	793621	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	30357	5.581	ng/uL	0.00
4) Pyridine-d5	3.819	84	436020	31.799	ng/ul	0.00
7) Phenol-d5	7.231	99	589190	35.186	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.378	67	408586	39.440	ng/ul	0.00
11) 2-Chlorophenol-d4	7.584	132	461959	36.547	ng/ul	0.00
15) 4-Methylphenol-d8	8.790	113	481749	36.418	ng/ul	0.00
21) Nitrobenzene-d5	9.242	128	214811	32.013	ng/ul	0.00
24) 2-Nitrophenol-d4	9.972	143	211316	26.983	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.536	165	395118	28.312	ng/ul	0.02
31) 4-Chloroaniline-d4	11.048	131	626126	34.926	ng/ul	0.01
46) Dimethylphthalate-d6	14.113	166	1408915	34.967	ng/ul	0.00
49) Acenaphthylene-d8	14.401	160	1612501	35.600	ng/ul	0.00
54) 4-Nitrophenol-d4	14.995	143	113743	21.392	ng/ul	0.04
60) Fluorene-d10	15.707	176	1186562	35.765	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.871	200	81026	12.408	ng/ul	0.01
73) Anthracene-d10	17.571	188	1779747	36.694	ng/ul	0.00
81) Pyrene-d10	19.795	212	1866819	38.547	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.936	264	1516474	37.880	ng/ul	0.00

Target Compounds	Qvalue

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

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