

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070623\
 Data File : BP016067.D
 Acq On : 07 Jul 2023 13:01
 Operator : MA/JU
 Sample : PB153590BL
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK590

Quant Time: Jul 07 20:25:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 06 23:18:26 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	224073	20.000	ng/u1	0.00
20) Naphthalene-d8	10.857	136	957184	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.681	164	627350	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	1459211	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1458297	20.000	ng/u1	0.02
88) Perylene-d12	24.092	264	1518083	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	39412	6.328	ng/uL	0.00
4) Pyridine-d5	3.799	84	479283	30.529	ng/u1	0.00
7) Phenol-d5	7.222	99	600705	31.332	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.352	67	426362	35.946	ng/u1	0.00
11) 2-Chlorophenol-d4	7.563	132	507096	35.039	ng/u1	0.00
15) 4-Methylphenol-d8	8.775	113	526363	34.753	ng/u1	0.00
21) Nitrobenzene-d5	9.228	128	244752	33.458	ng/u1	0.01
24) 2-Nitrophenol-d4	9.963	143	287444	33.668	ng/u1	0.02
28) 2,4-Dichlorophenol-d3	10.546	165	399281	26.244	ng/u1	0.04
31) 4-Chloroaniline-d4	11.051	131	634426	32.461	ng/u1	0.03
46) Dimethylphthalate-d6	14.098	166	1834790	37.839	ng/u1	0.01
49) Acenaphthylene-d8	14.375	160	1969850	36.138	ng/u1	0.00
54) 4-Nitrophenol-d4	15.016	143	151033	23.603	ng/u1	0.05
60) Fluorene-d10	15.686	176	1504310	37.677	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.875	200	215404	24.003	ng/u1	0.02
73) Anthracene-d10	17.551	188	2454612	36.826	ng/u1	0.00
81) Pyrene-d10	19.780	212	3115658	32.719	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.915	264	2786875	36.392	ng/u1	0.00

Target Compounds	Qvalue

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

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