

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015543.D
 Acq On : 15 Jun 2023 06:05
 Operator : MA/JU
 Sample : PB153339BL
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK339

Quant Time: Jun 15 06:36:12 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 15 05:36:46 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.087	152	100117	20.000	ng/u1	0.00
20) Naphthalene-d8	10.910	136	475805	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.728	164	336438	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.486	188	787991	20.000	ng/u1	0.00
79) Chrysene-d12	21.569	240	737969	20.000	ng/u1	0.00
88) Perylene-d12	24.115	264	757877	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	20877	7.722	ng/uL	0.00
4) Pyridine-d5	3.846	84	246630	35.119	ng/u1	0.00
7) Phenol-d5	7.240	99	352945	38.266	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	228366	41.458	ng/u1	0.00
11) 2-Chlorophenol-d4	7.610	132	277958	41.566	ng/u1	0.00
15) 4-Methylphenol-d8	8.799	113	309208	40.848	ng/u1	0.00
21) Nitrobenzene-d5	9.263	128	152838	40.915	ng/u1	0.00
24) 2-Nitrophenol-d4	9.993	143	178501	41.843	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.534	165	308955	39.702	ng/u1	0.00
31) 4-Chloroaniline-d4	11.057	131	359876	32.218	ng/u1	0.00
46) Dimethylphthalate-d6	14.134	166	1133212	42.714	ng/u1	0.00
49) Acenaphthylene-d8	14.422	160	1217004	41.952	ng/u1	0.00
54) 4-Nitrophenol-d4	14.934	143	128931	27.138	ng/u1	0.00
60) Fluorene-d10	15.722	176	936987	41.882	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.851	200	172951	34.646	ng/u1	0.00
73) Anthracene-d10	17.586	188	1539437	42.938	ng/u1	0.00
81) Pyrene-d10	19.810	212	1891086	47.879	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.951	264	1649218	43.370	ng/u1	0.00

Target Compounds Qvalue

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

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