

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061223\
 Data File : BP015478.D
 Acq On : 12 Jun 2023 18:37
 Operator : MA/JU
 Sample : 03062-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZEL3

Quant Time: Jun 12 23:23:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 09 23:33:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.093	152	133581	20.000	ng/ul	0.00
20) Naphthalene-d8	10.922	136	622552	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	437473	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.492	188	1033131	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240	1015725	20.000	ng/ul	0.00
88) Perylene-d12	24.116	264	1064470	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	11731	3.252	ng/uL	0.00
4) Pyridine-d5	3.852	84	137128	14.635	ng/ul	0.00
7) Phenol-d5	7.246	99	212441	17.263	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	136752	18.607	ng/ul	0.00
11) 2-Chlorophenol-d4	7.622	132	167152	18.734	ng/ul	0.00
15) 4-Methylphenol-d8	8.805	113	182783	18.097	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128	87331	17.868	ng/ul	0.00
24) 2-Nitrophenol-d4	9.999	143	98785	17.698	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.546	165	170695	16.765	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	258581	17.693	ng/ul	0.00
46) Dimethylphthalate-d6	14.140	166	678758	19.676	ng/ul	0.00
49) Acenaphthylene-d8	14.434	160	733176	19.437	ng/ul	0.00
54) 4-Nitrophenol-d4	14.957	143	68271	11.051	ng/ul	0.02
60) Fluorene-d10	15.728	176	577480	19.851	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	78650	12.017	ng/ul	0.00
73) Anthracene-d10	17.592	188	940340	20.005	ng/ul	0.00
81) Pyrene-d10	19.810	212	1172468	21.567	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	1059584	19.838	ng/ul	0.00

Target Compounds Qvalue

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

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