

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060923\
 Data File : BP015466.D
 Acq On : 10 Jun 2023 00:11
 Operator : MA/JU
 Sample : 03097-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZEN3

Quant Time: Jun 10 00:41:38 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	137532	20.000	ng/u1	0.00
20) Naphthalene-d8	10.922	136	641086	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.734	164	453558	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.492	188	1063717	20.000	ng/u1	0.00
79) Chrysene-d12	21.575	240	1051801	20.000	ng/u1	0.00
88) Perylene-d12	24.121	264	1082382	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.411	96	15261	4.109	ng/uL	0.00
4) Pyridine-d5	3.852	84	148850	15.429	ng/u1	0.00
7) Phenol-d5	7.246	99	154741	12.213	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	469048	61.986	ng/u1	0.00
11) 2-Chlorophenol-d4	7.616	132	430460	46.859	ng/u1	0.00
15) 4-Methylphenol-d8	8.805	113	314816	30.274	ng/u1	0.00
21) Nitrobenzene-d5	9.275	128	316537	62.890	ng/u1	0.00
24) 2-Nitrophenol-d4	9.999	143	349851	60.867	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	552917	52.734	ng/u1	0.00
31) 4-Chloroaniline-d4	11.063	131	727224	48.321	ng/u1	0.00
46) Dimethylphthalate-d6	14.146	166	2408547	67.342	ng/u1	0.00
49) Acenaphthylene-d8	14.434	160	2526005	64.590	ng/u1	0.00
54) 4-Nitrophenol-d4	14.945	143	63014	9.838	ng/u1	0.00
60) Fluorene-d10	15.734	176	2014809	66.803	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	441870	65.572	ng/u1	0.00
73) Anthracene-d10	17.592	188	3491817	72.148	ng/u1	0.00
81) Pyrene-d10	19.816	212	4210801	74.800	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.963	264	4080034	75.126	ng/u1	0.00

Target Compounds Qvalue

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

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