

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092123\
 Data File : BP017273.D
 Acq On : 21 Sep 2023 18:21
 Operator : MA/JU
 Sample : 04387-13
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AH1

Quant Time: Sep 21 23:58:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 21 23:51:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	168058	20.000	ng/u1	0.00
20) Naphthalene-d8	10.763	136	665152	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	388281	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.369	188	872678	20.000	ng/u1	0.00
79) Chrysene-d12	21.474	240	873706	20.000	ng/u1	0.00
88) Perylene-d12	23.998	264	999717	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	34024	8.316	ng/uL	0.00
4) Pyridine-d5	3.740	84	574625	50.236	ng/u1	0.00
7) Phenol-d5	7.098	99	794263	57.545	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	448993	53.902	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	610620	55.779	ng/u1	0.00
15) 4-Methylphenol-d8	8.657	113	658455	61.361	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	323784	67.334	ng/u1	0.00
24) 2-Nitrophenol-d4	9.851	143	387103	74.139	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	682859	69.223	ng/u1	0.00
31) 4-Chloroaniline-d4	10.928	131	962140	67.331	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	2306252	82.913	ng/u1	0.00
49) Acenaphthylene-d8	14.298	160	2549649	79.699	ng/u1	0.00
54) 4-Nitrophenol-d4	14.833	143	414156	85.465	ng/u1	0.00
60) Fluorene-d10	15.598	176	1998773	83.469	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	333862	67.878	ng/u1	0.00
73) Anthracene-d10	17.474	188	3134279	80.981	ng/u1	0.00
81) Pyrene-d10	19.715	212	3927744	83.131	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	4105552	85.295	ng/u1	0.00

Target Compounds	Qvalue

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

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