

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
 Data File : BP017283.D
 Acq On : 22 Sep 2023 10:38
 Operator : MA/JU
 Sample : 04387-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AF3

Quant Time: Sep 22 12:44:27 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	7.934	152	218719	20.000	ng/u1	0.00
20) Naphthalene-d8	10.757	136	881481	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	507523	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.368	188	1164427	20.000	ng/u1	0.00
79) Chrysene-d12	21.474	240	1172216	20.000	ng/u1	0.00
88) Perylene-d12	23.992	264	1327215	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.304	96	20001	3.756	ng/uL	0.00
4) Pyridine-d5	3.740	84	304010	20.422	ng/u1	0.00
7) Phenol-d5	7.093	99	403802	22.479	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	237153	21.876	ng/u1	0.00
11) 2-Chlorophenol-d4	7.457	132	315461	22.142	ng/u1	0.00
15) 4-Methylphenol-d8	8.651	113	327784	23.471	ng/u1	0.00
21) Nitrobenzene-d5	9.134	128	154388	24.227	ng/u1	0.00
24) 2-Nitrophenol-d4	9.845	143	173463	25.069	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.375	165	321961	24.628	ng/u1	0.00
31) 4-Chloroaniline-d4	10.922	131	437159	23.085	ng/u1	0.00
46) Dimethylphthalate-d6	14.010	166	990767	27.251	ng/u1	0.00
49) Acenaphthylene-d8	14.292	160	1114479	26.652	ng/u1	0.00
54) 4-Nitrophenol-d4	14.827	143	142481	22.494	ng/u1	0.00
60) Fluorene-d10	15.592	176	865646	27.656	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200	108764	16.572	ng/u1	0.00
73) Anthracene-d10	17.468	188	1390117	26.918	ng/u1	0.00
81) Pyrene-d10	19.709	212	1733960	27.354	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.821	264	1747942	27.354	ng/u1	-0.01

Target Compounds Qvalue

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

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