

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017177.D
 Acq On : 14 Sep 2023 15:13
 Operator : MA/JU
 Sample : 04227-19
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJ99

Quant Time: Sep 14 22:48:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 12 23:27:27 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	172687	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.775	136	680195	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.610	164	396328	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	879305	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	837549	20.000	ng/u1	0.00
88) Perylene-d12	24.016	264	961340	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	21722	5.286	ng/uL	0.00
4) Pyridine-d5	3.752	84	82120	6.986	ng/u1	0.00
7) Phenol-d5	7.099	99	103389	7.363	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	293259	33.414	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.469	132	291003	25.978	ng/u1	0.00
15) 4-Methylphenol-d8	8.658	113	185237	15.918	ng/u1	-0.01
21) Nitrobenzene-d5	9.146	128	171941	36.100	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.863	143	176329	34.678	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	299302	30.436	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	341977	23.072	ng/u1	0.00
46) Dimethylphthalate-d6	14.028	166	1026147	36.384	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	1150042	35.936	ng/u1	0.00
54) 4-Nitrophenol-d4	14.840	143	23901	4.385	ng/u1	0.00
60) Fluorene-d10	15.610	176	907777	37.225	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.740	200	127668	25.987	ng/u1	0.00
73) Anthracene-d10	17.481	188	1465989	39.240	ng/u1	0.00
81) Pyrene-d10	19.722	212	1750773	39.858	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.851	264	1790848	39.469	ng/u1	0.00

Target Compounds	Qvalue

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

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