

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017153.D
 Acq On : 13 Sep 2023 17:25
 Operator : MA/JU
 Sample : 04324-09
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHE9

Quant Time: Sep 13 23:19:43 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	7.952	152	243148	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	999439	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.610	164	623898	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	1408033	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	1158661	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	1077947	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	23002	3.975	ng/uL	0.00
4) Pyridine-d5	3.752	84	221845	13.403	ng/u1	0.00
7) Phenol-d5	7.105	99	168174	8.506	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	433149	35.052	ng/u1	0.00
11) 2-Chlorophenol-d4	7.475	132	432576	27.426	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	298691	18.229	ng/u1	0.00
21) Nitrobenzene-d5	9.152	128	257761	36.832	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	273501	36.607	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	475758	32.927	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	391322	17.968	ng/u1	0.00
46) Dimethylphthalate-d6	14.028	166	1748301	39.378	ng/u1	0.00
49) Acenaphthylene-d8	14.310	160	1862047	36.962	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	54253	6.323	ng/u1	0.00
60) Fluorene-d10	15.610	176	1510818	39.356	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.734	200	263983	33.557	ng/u1	-0.01
73) Anthracene-d10	17.481	188	2584524	43.202	ng/u1	0.00
81) Pyrene-d10	19.722	212	3116804	51.292	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	2356221	46.312	ng/u1	-0.01

Target Compounds Qvalue

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

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