

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017182.D
 Acq On : 14 Sep 2023 18:05
 Operator : MA/JU
 Sample : 04324-10
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHR6

Quant Time: Sep 14 23:57:42 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	231914	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	933832	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.610	164	534730	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.381	188	1160687	20.000	ng/u1	0.00
79) Chrysene-d12	21.480	240	1056057	20.000	ng/u1	0.00
88) Perylene-d12	24.004	264	1198467	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	31826	5.767	ng/uL	0.00
4) Pyridine-d5	3.746	84	149560	9.473	ng/u1	0.00
7) Phenol-d5	7.099	99	113800	6.035	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	356846	30.276	ng/u1	0.00
11) 2-Chlorophenol-d4	7.469	132	337196	22.414	ng/u1	0.00
15) 4-Methylphenol-d8	8.658	113	215196	13.769	ng/u1	-0.01
21) Nitrobenzene-d5	9.146	128	209653	32.062	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.863	143	217181	31.111	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	382131	28.305	ng/u1	0.00
31) 4-Chloroaniline-d4	10.934	131	273134	13.423	ng/u1	-0.01
46) Dimethylphthalate-d6	14.022	166	1333444	35.042	ng/u1	-0.01
49) Acenaphthylene-d8	14.310	160	1409356	32.641	ng/u1	0.00
54) 4-Nitrophenol-d4	14.834	143	30857	4.196	ng/u1	0.00
60) Fluorene-d10	15.604	176	1145399	34.812	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.734	200	189478	29.219	ng/u1	-0.01
73) Anthracene-d10	17.481	188	2012547	40.810	ng/u1	0.00
81) Pyrene-d10	19.722	212	2421689	43.725	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.839	264	2386993	42.199	ng/u1	-0.01

Target Compounds Qvalue

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

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