

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017614.D
 Acq On : 06 Oct 2023 22:12
 Operator : MA/JU
 Sample : 04624-18
 Misc :
 ALS Vial : 94 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHB1

Quant Time: Oct 06 23:16:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 06 05:55:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	83198	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	335447	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.622	164	193068	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	463526	20.000	ng/u1	0.00
79) Chrysene-d12	21.551	240	492815	20.000	ng/u1	0.00
88) Perylene-d12	24.127	264	546986	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	13318	6.400	ng/uL	0.00
4) Pyridine-d5	3.722	84	86371	15.856	ng/u1	0.00
7) Phenol-d5	7.093	99	73855	10.551	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	204919	47.975	ng/u1	0.00
11) 2-Chlorophenol-d4	7.452	132	209325	37.978	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	133190	23.767	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	126541	49.034	ng/u1	0.00
24) 2-Nitrophenol-d4	9.857	143	137655	47.293	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	232744	43.132	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	239203	31.118	ng/u1	0.00
46) Dimethylphthalate-d6	14.040	166	736693	48.918	ng/u1	0.00
49) Acenaphthylene-d8	14.322	160	818540	49.291	ng/u1	0.00
54) 4-Nitrophenol-d4	14.881	143	18912	7.171	ng/u1	0.01
60) Fluorene-d10	15.628	176	656759	51.414	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	101718	35.200	ng/u1	0.00
73) Anthracene-d10	17.516	188	1098968	51.219	ng/u1	0.00
81) Pyrene-d10	19.769	212	1395777	50.970	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.957	264	1459391	52.868	ng/u1	0.00

Target Compounds Qvalue

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

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