

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017608.D
 Acq On : 06 Oct 2023 18:46
 Operator : MA/JU
 Sample : 04624-16
 Misc :
 ALS Vial : 88 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHA5

Quant Time: Oct 06 23:15:15 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	7.928	152	91184	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	373152	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	225712	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.416	188	545746	20.000	ng/u1	0.00
79) Chrysene-d12	21.551	240	554123	20.000	ng/u1	0.00
88) Perylene-d12	24.127	264	606452	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	11119	4.875	ng/uL	0.00
4) Pyridine-d5	3.728	84	80901	13.551	ng/u1	0.00
7) Phenol-d5	7.093	99	68659	8.949	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	182032	38.884	ng/u1	0.00
11) 2-Chlorophenol-d4	7.451	132	186939	30.946	ng/u1	0.00
15) 4-Methylphenol-d8	8.651	113	120337	19.593	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	114099	39.745	ng/u1	0.00
24) 2-Nitrophenol-d4	9.857	143	117381	36.253	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	210135	35.007	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	228506	26.723	ng/u1	0.00
46) Dimethylphthalate-d6	14.039	166	710817	40.373	ng/u1	0.00
49) Acenaphthylene-d8	14.322	160	765003	39.404	ng/u1	0.00
54) 4-Nitrophenol-d4	14.886	143	19312	6.264	ng/u1	0.02
60) Fluorene-d10	15.627	176	625099	41.858	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	97764	28.734	ng/u1	0.00
73) Anthracene-d10	17.516	188	1094320	43.319	ng/u1	0.00
81) Pyrene-d10	19.768	212	1400683	45.490	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.956	264	1401979	45.808	ng/u1	0.00

Target Compounds Qvalue

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

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