

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017616.D
 Acq On : 06 Oct 2023 23:21
 Operator : MA/JU
 Sample : 04624-20
 Misc :
 ALS Vial : 96 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBK86

Quant Time: Oct 07 00:11:12 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	84059	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	346612	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.622	164	219843	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.410	188	507996	20.000	ng/u1	0.00
79) Chrysene-d12	21.545	240	515322	20.000	ng/u1	0.00
88) Perylene-d12	24.121	264	569726	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	10241	4.871	ng/uL	0.00
4) Pyridine-d5	3.728	84	40840	7.421	ng/u1	0.00
7) Phenol-d5	7.093	99	59883	8.467	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	164459	38.108	ng/u1	0.00
11) 2-Chlorophenol-d4	7.452	132	168173	30.199	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	113772	20.094	ng/u1	0.00
21) Nitrobenzene-d5	9.140	128	109177	40.943	ng/u1	0.00
24) 2-Nitrophenol-d4	9.857	143	95957	31.905	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	189089	33.913	ng/u1	0.00
31) 4-Chloroaniline-d4	10.940	131	181468	22.847	ng/u1	0.00
46) Dimethylphthalate-d6	14.040	166	680067	39.658	ng/u1	0.00
49) Acenaphthylene-d8	14.322	160	745461	39.423	ng/u1	0.00
54) 4-Nitrophenol-d4	14.892	143	13733	4.573	ng/u1	0.02
60) Fluorene-d10	15.628	176	613822	42.201	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.769	200	68350	21.582	ng/u1	0.00
73) Anthracene-d10	17.510	188	1126884	47.923	ng/u1	0.00
81) Pyrene-d10	19.769	212	1384781	48.360	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.951	264	1403804	48.825	ng/u1	0.00

Target Compounds	Qvalue

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

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