

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017563.D
 Acq On : 05 Oct 2023 13:40
 Operator : MA/JU
 Sample : 04679-04
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PMW-4(20230927)

Quant Time: Oct 06 02:34:11 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 04 16:33:47 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	109726	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	442308	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.628	164	278527	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.422	188	664169	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	689349	20.000	ng/u1	0.00
88) Perylene-d12	24.139	264	773349	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	11258	4.102	ng/uL	0.00
4) Pyridine-d5	3.728	84	104039	14.482	ng/u1	0.00
7) Phenol-d5	7.099	99	83838	9.081	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	219351	38.938	ng/u1	0.00
11) 2-Chlorophenol-d4	7.457	132	222522	30.612	ng/u1	0.00
15) 4-Methylphenol-d8	8.657	113	176700	23.908	ng/u1	-0.01
21) Nitrobenzene-d5	9.146	128	138621	40.737	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	151575	39.494	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	276239	38.824	ng/u1	0.00
31) 4-Chloroaniline-d4	10.945	131	313765	30.956	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	969277	44.614	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	1022586	42.684	ng/u1	0.00
54) 4-Nitrophenol-d4	14.881	143	25258	6.639	ng/u1	0.00
60) Fluorene-d10	15.633	176	823351	44.679	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	142160	34.333	ng/u1	0.00
73) Anthracene-d10	17.522	188	1349413	43.892	ng/u1	0.00
81) Pyrene-d10	19.774	212	1775534	46.353	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	1778998	45.583	ng/u1	0.00

Target Compounds	Qvalue

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

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