

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017566.D
 Acq On : 05 Oct 2023 15:22
 Operator : MA/JU
 Sample : 04679-17
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PMW-9D(20230928)

Quant Time: Oct 06 02:42:43 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	94516	20.000	ng/u1	0.00
20) Naphthalene-d8	10.781	136	367313	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	239688	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.422	188	557033	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	574273	20.000	ng/u1	0.00
88) Perylene-d12	24.133	264	646427	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	7975	3.373	ng/uL	0.00
4) Pyridine-d5	3.728	84	62362	10.078	ng/u1	0.00
7) Phenol-d5	7.099	99	57986	7.292	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	174153	35.890	ng/u1	0.00
11) 2-Chlorophenol-d4	7.458	132	175951	28.100	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	115045	18.071	ng/u1	0.00
21) Nitrobenzene-d5	9.146	128	109322	38.687	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	117860	36.979	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	205435	34.768	ng/u1	0.00
31) 4-Chloroaniline-d4	10.946	131	252901	30.046	ng/u1	0.00
46) Dimethylphthalate-d6	14.045	166	741382	39.654	ng/u1	0.00
49) Acenaphthylene-d8	14.328	160	785605	38.106	ng/u1	0.00
54) 4-Nitrophenol-d4	14.887	143	20494	6.259	ng/u1	0.01
60) Fluorene-d10	15.634	176	634859	40.033	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	101534	29.238	ng/u1	0.00
73) Anthracene-d10	17.522	188	1072190	41.583	ng/u1	0.00
81) Pyrene-d10	19.775	212	1369644	42.921	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	1397508	42.838	ng/u1	0.00

Target Compounds	Qvalue

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

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