

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017550.D
 Acq On : 05 Oct 2023 03:56
 Operator : MA/JU
 Sample : 04679-16
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

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

Quant Time: Oct 05 04:32:33 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	7.946	152	116130	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	481080	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.634	164	313508	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.422	188	733350	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	748179	20.000	ng/u1	0.00
88) Perylene-d12	24.139	264	833750	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	13955	4.804	ng/uL	0.00
4) Pyridine-d5	3.728	84	122182	16.070	ng/u1	0.00
7) Phenol-d5	7.099	99	134482	13.764	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	277576	46.557	ng/u1	0.00
11) 2-Chlorophenol-d4	7.463	132	287103	37.318	ng/u1	0.00
15) 4-Methylphenol-d8	8.663	113	261419	33.421	ng/u1	0.00
21) Nitrobenzene-d5	9.146	128	180377	48.736	ng/u1	0.00
24) 2-Nitrophenol-d4	9.863	143	199620	47.821	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	359405	46.442	ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	493595	44.774	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	1420736	58.097	ng/u1	0.00
49) Acenaphthylene-d8	14.334	160	1411687	52.351	ng/u1	0.00
54) 4-Nitrophenol-d4	14.887	143	48324	11.284	ng/u1	0.01
60) Fluorene-d10	15.639	176	1139451	54.933	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	201470	44.067	ng/u1	0.00
73) Anthracene-d10	17.522	188	1997812	58.853	ng/u1	0.00
81) Pyrene-d10	19.780	212	2483589	59.739	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	2516858	59.817	ng/u1	0.00

Target Compounds Qvalue

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

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