

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017558.D
 Acq On : 05 Oct 2023 10:50
 Operator : MA/JU
 Sample : 04679-22
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EB(20230928)

Quant Time: Oct 06 01:58:18 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	102515	20.000	ng/ul	0.00
20) Naphthalene-d8	10.781	136	411221	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.640	164	259220	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.428	188	587689	20.000	ng/ul	0.00
79) Chrysene-d12	21.563	240	611393	20.000	ng/ul	0.00
88) Perylene-d12	24.145	264	694467	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	9037	3.524	ng/uL	0.00
4) Pyridine-d5	3.734	84	36042	5.370	ng/ul	0.00
7) Phenol-d5	7.105	99	46863	5.433	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	149281	28.364	ng/ul	0.00
11) 2-Chlorophenol-d4	7.463	132	144914	21.338	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	89494	12.961	ng/ul	0.00
21) Nitrobenzene-d5	9.152	128	91671	28.977	ng/ul	0.00
24) 2-Nitrophenol-d4	9.869	143	95867	26.867	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.399	165	163923	24.780	ng/ul	0.00
31) 4-Chloroaniline-d4	10.951	131	235170	24.956	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	599789	29.663	ng/ul	0.00
49) Acenaphthylene-d8	14.334	160	644317	28.898	ng/ul	0.00
54) 4-Nitrophenol-d4	14.892	143	14698	4.151	ng/ul	0.02
60) Fluorene-d10	15.639	176	523482	30.523	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	75352	20.567	ng/ul	0.00
73) Anthracene-d10	17.528	188	894733	32.890	ng/ul	0.00
81) Pyrene-d10	19.780	212	1150766	33.873	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.974	264	1175327	33.536	ng/ul	0.00

Target Compounds Qvalue

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

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