

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017576.D
 Acq On : 05 Oct 2023 21:05
 Operator : MA/JU
 Sample : 04607-01
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Oct 06 02:03:34 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	7.934	152	77141	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.769	136	307547	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.628	164	197385	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.416	188	462126	20.000	ng/ul	-0.01
79) Chrysene-d12	21.551	240	475922	20.000	ng/ul	-0.01
88) Perylene-d12	24.127	264	531035	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	17939	9.297	ng/uL	0.00
4) Pyridine-d5	3.728	84	53658	10.624	ng/ul	0.00
7) Phenol-d5	7.099	99	50116	7.722	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	162203	40.956	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.458	132	156148	30.554	ng/ul	0.00
15) 4-Methylphenol-d8	8.658	113	97579	18.780	ng/ul	-0.01
21) Nitrobenzene-d5	9.146	128	99407	42.014	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	102859	38.544	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.393	165	182411	36.871	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	234931	33.335	ng/ul	0.00
46) Dimethylphthalate-d6	14.046	166	768317	49.902	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	755601	44.506	ng/ul	0.00
54) 4-Nitrophenol-d4	14.887	143	17592	6.525	ng/ul	0.01
60) Fluorene-d10	15.634	176	665988	50.997	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.775	200	105253	36.533	ng/ul	0.00
73) Anthracene-d10	17.516	188	1177801	55.060	ng/ul	-0.01
81) Pyrene-d10	19.775	212	1479938	55.962	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.963	264	1510881	56.377	ng/ul	-0.01

Target Compounds	Qvalue

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

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