

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017573.D
 Acq On : 05 Oct 2023 19:22
 Operator : MA/JU
 Sample : 04588-17
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Oct 06 02:03:00 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	79593	20.000	ng/ul	0.00
20) Naphthalene-d8	10.775	136	319590	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.628	164	201671	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.416	188	472825	20.000	ng/ul	-0.01
79) Chrysene-d12	21.557	240	482051	20.000	ng/ul	0.00
88) Perylene-d12	24.133	264	533463	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	17981	9.032	ng/uL	0.00
4) Pyridine-d5	3.734	84	59533	11.424	ng/ul	0.00
7) Phenol-d5	7.099	99	49003	7.317	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	149025	36.469	ng/ul	0.00
11) 2-Chlorophenol-d4	7.457	132	149716	28.393	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	92045	17.169	ng/ul	0.00
21) Nitrobenzene-d5	9.146	128	86433	35.154	ng/ul	0.00
24) 2-Nitrophenol-d4	9.863	143	87658	31.610	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.393	165	167139	32.511	ng/ul	0.00
31) 4-Chloroaniline-d4	10.946	131	219939	30.032	ng/ul	0.00
46) Dimethylphthalate-d6	14.045	166	592408	37.659	ng/ul	0.00
49) Acenaphthylene-d8	14.328	160	640220	36.908	ng/ul	0.00
54) 4-Nitrophenol-d4	14.892	143	12653	4.593	ng/ul	0.02
60) Fluorene-d10	15.634	176	518651	38.871	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.781	200	65783	22.317	ng/ul	0.00
73) Anthracene-d10	17.522	188	856608	39.138	ng/ul	0.00
81) Pyrene-d10	19.774	212	1079753	40.310	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.962	264	1101176	40.903	ng/ul	-0.01

Target Compounds Qvalue

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

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