

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017582.D
 Acq On : 06 Oct 2023 00:29
 Operator : MA/JU
 Sample : 04588-21
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Oct 06 02:04:44 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	83787	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.769	136	349547	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.628	164	220263	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.416	188	523585	20.000	ng/ul	-0.01
79) Chrysene-d12	21.551	240	544625	20.000	ng/ul	-0.01
88) Perylene-d12	24.121	264	611060	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	12077	5.763	ng/uL	0.00
4) Pyridine-d5	3.728	84	79100	14.419	ng/ul	0.00
7) Phenol-d5	7.093	99	60771	8.621	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67	197253	45.855	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.452	132	189060	34.060	ng/ul	-0.01
15) 4-Methylphenol-d8	8.658	113	111701	19.793	ng/ul	-0.01
21) Nitrobenzene-d5	9.146	128	117837	43.819	ng/ul	0.00
24) 2-Nitrophenol-d4	9.858	143	127030	41.882	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.387	165	214676	38.179	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.940	131	264107	32.972	ng/ul	-0.01
46) Dimethylphthalate-d6	14.046	166	819960	47.724	ng/ul	0.00
49) Acenaphthylene-d8	14.322	160	872353	46.045	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.887	143	17106	5.685	ng/ul	0.01
60) Fluorene-d10	15.628	176	718497	49.303	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.775	200	107061	32.799	ng/ul	0.00
73) Anthracene-d10	17.516	188	1190866	49.136	ng/ul	-0.01
81) Pyrene-d10	19.769	212	1521262	50.268	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.957	264	1508732	48.925	ng/ul	-0.02

Target Compounds Qvalue

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

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