

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018156.D
 Acq On : 14 Nov 2023 15:36
 Operator : MA/JU
 Sample : 05337-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 15 00:38:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	85973	20.000	ng/ul	0.00
20) Naphthalene-d8	10.605	136	354658	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	223398	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188	514933	20.000	ng/ul	-0.01
79) Chrysene-d12	21.380	240	539895	20.000	ng/ul	-0.01
88) Perylene-d12	23.851	264	605329	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	7241	2.971	ng/uL	0.00
4) Pyridine-d5	3.623	84	13407	2.110	ng/ul	0.02
7) Phenol-d5	6.964	99	39183	4.721	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	137223	28.133	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	126043	19.176	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	79340	11.665	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	82254	25.742	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	88006	24.311	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	138235	21.426	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	148405	16.773	ng/ul	0.00
46) Dimethylphthalate-d6	13.893	166	611692	29.676	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	603789	26.894	ng/ul	0.00
54) 4-Nitrophenol-d4	14.781	143	6468	2.100	ng/ul	0.04
60) Fluorene-d10	15.469	176	494210	29.041	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	90942	24.430	ng/ul	-0.04
73) Anthracene-d10	17.351	188	872311	31.322	ng/ul	0.00
81) Pyrene-d10	19.604	212	1158169	32.115	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1176366	33.421	ng/ul	-0.02

Target Compounds Qvalue

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

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