

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018163.D
 Acq On : 14 Nov 2023 19:32
 Operator : MA/JU
 Sample : 05337-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDJ4

Quant Time: Nov 15 01:33:33 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	77019	20.000	ng/u1	0.00
20) Naphthalene-d8	10.604	136	316275	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.463	164	209866	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.245	188	497858	20.000	ng/u1	-0.01
79) Chrysene-d12	21.374	240	515263	20.000	ng/u1	-0.02
88) Perylene-d12	23.845	264	573629	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.163	96	6491	2.973	ng/uL	0.00
4) Pyridine-d5	3.605	84	59880	10.519	ng/u1	0.00
7) Phenol-d5	7.016	99	2394	0.322	ng/u1	0.06
9) Bis-(2-Chloroethyl)eth...	7.140	67	122878	28.120	ng/u1	0.00
11) 2-Chlorophenol-d4	0.000	132	0d	0.000	ng/u1	
15) 4-Methylphenol-d8	8.534	113	17633	2.894	ng/u1	0.02
21) Nitrobenzene-d5	8.992	128	74110	26.008	ng/u1	0.00
24) 2-Nitrophenol-d4	0.000	143	0d	0.000	ng/u1	
28) 2,4-Dichlorophenol-d3	0.000	165	0d	0.000	ng/u1	
31) 4-Chloroaniline-d4	10.787	131	186617	23.651	ng/u1	0.00
46) Dimethylphthalate-d6	13.886	166	594792	30.717	ng/u1	-0.01
49) Acenaphthylene-d8	14.163	160	590195	27.984	ng/u1	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/u1	
60) Fluorene-d10	15.469	176	507181	31.725	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/u1	
73) Anthracene-d10	17.351	188	897827	33.344	ng/u1	0.00
81) Pyrene-d10	19.604	212	1170225	34.001	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1188333	35.627	ng/u1	-0.02
Target Compounds						Qvalue
2) 1,4-Dioxane	3.199	88	3226	1.354	ng/uL	90

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

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