

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020725\
 Data File : BP024003.D
 Acq On : 07 Feb 2025 15:10
 Operator : RC/JU
 Sample : Q1229-18
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29C4

Quant Time: Feb 12 13:14:22 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 03 11:19:36 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	376165	20.000	ng/u1	0.00
20) Naphthalene-d8	10.534	136	1546851	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.387	164	935855	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.181	188	1931594	20.000	ng/u1	0.00
79) Chrysene-d12	21.627	240	1898892	20.000	ng/u1	-0.01
88) Perylene-d12	24.986	264	1935169	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	42742	4.721	ng/uL	0.00
4) Pyridine-d5	3.705	84	276119	10.569	ng/u1	0.00
7) Phenol-d5	6.952	99	252145	7.574	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	554272	31.456	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	684897	26.200	ng/u1	0.00
15) 4-Methylphenol-d8	8.469	113	521378	19.173	ng/u1	0.00
21) Nitrobenzene-d5	8.910	128	394286	31.859	ng/u1	0.00
24) 2-Nitrophenol-d4	9.628	143	398128	29.814	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	728713	29.746	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	1046446	28.563	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2440130	34.956	ng/u1	0.00
49) Acenaphthylene-d8	14.075	160	2790856	35.448	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	64361	4.643	ng/u1	0.00
60) Fluorene-d10	15.392	176	2245148	38.193	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	292385	24.536	ng/u1	0.00
73) Anthracene-d10	17.281	188	3762261	39.669	ng/u1	0.00
81) Pyrene-d10	19.651	212	4398446	43.230	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.751	264	4087586	40.484	ng/u1	-0.02

Target Compounds	Qvalue

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

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