

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
 Data File : BP023919.D
 Acq On : 04 Feb 2025 19:28
 Operator : RC/JU
 Sample : Q1202-15
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29A7

Quant Time: Feb 05 07:35:05 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	7.775	152	478563	20.000	ng/u1	0.00
20) Naphthalene-d8	10.546	136	2042356	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.392	164	1272890	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.181	188	2507860	20.000	ng/u1	0.00
79) Chrysene-d12	21.616	240	2428643	20.000	ng/u1	-0.02
88) Perylene-d12	24.986	264	2591585	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	56195	4.879	ng/uL	0.00
4) Pyridine-d5	3.705	84	308922	9.295	ng/u1	0.00
7) Phenol-d5	6.958	99	379562	8.962	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	710703	31.703	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	956168	28.750	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	718622	20.772	ng/u1	0.00
21) Nitrobenzene-d5	8.916	128	520643	31.863	ng/u1	0.00
24) 2-Nitrophenol-d4	9.634	143	546445	30.993	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.175	165	1010043	31.227	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	1287550	26.618	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	3278489	34.530	ng/u1	0.00
49) Acenaphthylene-d8	14.081	160	3667540	34.249	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	141436	7.501	ng/u1	0.00
60) Fluorene-d10	15.398	176	2901786	36.293	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	395502	25.563	ng/u1	0.00
73) Anthracene-d10	17.281	188	4649460	37.759	ng/u1	0.00
81) Pyrene-d10	19.645	212	5374383	41.300	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.757	264	5274303	39.006	ng/u1	-0.02

Target Compounds Qvalue

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

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