

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
 Data File : BP023911.D
 Acq On : 04 Feb 2025 13:56
 Operator : RC/JU
 Sample : Q1202-11
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29A2

Quant Time: Feb 05 07:33:40 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	7.746	152	557244	20.000	ng/u1	-0.03
20) Naphthalene-d8	10.522	136	2427457	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.392	164	1555220	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.204	188	2911233	20.000	ng/u1	0.02
79) Chrysene-d12	21.651	240	2847629	20.000	ng/u1	0.01
88) Perylene-d12	25.027	264	3010940	20.000	ng/u1	0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	72023	5.370	ng/uL	-0.02
4) Pyridine-d5	3.687	84	297692	7.692	ng/u1	-0.02
7) Phenol-d5	6.922	99	416222	8.440	ng/u1	-0.04
9) Bis-(2-Chloroethyl)eth...	7.075	67	817570	31.321	ng/u1	-0.03
11) 2-Chlorophenol-d4	7.287	132	1109104	28.640	ng/u1	-0.03
15) 4-Methylphenol-d8	8.451	113	792658	19.676	ng/u1	-0.02
21) Nitrobenzene-d5	8.893	128	608479	31.331	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.604	143	669061	31.927	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.146	165	1172115	30.489	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.657	131	1486227	25.851	ng/u1	-0.02
46) Dimethylphthalate-d6	13.798	166	3764272	32.449	ng/u1	0.00
49) Acenaphthylene-d8	14.086	160	4179019	31.941	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	180995	7.856	ng/u1	0.00
60) Fluorene-d10	15.398	176	3283065	33.607	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.528	200	555760	30.944	ng/u1	0.02
73) Anthracene-d10	17.310	188	5422074	37.932	ng/u1	0.02
81) Pyrene-d10	19.686	212	6176969	40.483	ng/u1	0.02
92) Benzo(a)pyrene-d12	24.804	264	6161199	39.219	ng/u1	0.03

Target Compounds	Qvalue

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

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