

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
 Data File : BP023923.D
 Acq On : 04 Feb 2025 22:10
 Operator : RC/JU
 Sample : Q1247-03
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A6328

Quant Time: Feb 05 07:36:02 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	485905	20.000	ng/u1	0.00
20) Naphthalene-d8	10.551	136	2117933	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.392	164	1339936	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.175	188	2560561	20.000	ng/u1	-0.01
79) Chrysene-d12	21.627	240	2426586	20.000	ng/u1	-0.01
88) Perylene-d12	24.986	264	2535590	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	52830	4.517	ng/uL	0.00
4) Pyridine-d5	3.711	84	125620	3.723	ng/u1	0.00
7) Phenol-d5	6.957	99	307185	7.144	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	693888	30.486	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	912344	27.018	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	626347	17.831	ng/u1	0.00
21) Nitrobenzene-d5	8.916	128	507683	29.961	ng/u1	0.00
24) 2-Nitrophenol-d4	9.634	143	535957	29.313	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.175	165	1011174	30.147	ng/u1	0.00
31) 4-Chloroaniline-d4	10.681	131	615512	12.270	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	3252583	32.543	ng/u1	0.00
49) Acenaphthylene-d8	14.081	160	3634349	32.241	ng/u1	0.00
54) 4-Nitrophenol-d4	14.604	143	117403	5.915	ng/u1	0.00
60) Fluorene-d10	15.392	176	2867311	34.067	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	374060	23.680	ng/u1	0.00
73) Anthracene-d10	17.280	188	4175686	33.213	ng/u1	0.00
81) Pyrene-d10	19.657	212	5089440	39.143	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.762	264	4846032	36.630	ng/u1	-0.01

Target Compounds	Qvalue

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

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