

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013125\
 Data File : BP023857.D
 Acq On : 01 Feb 2025 14:08
 Operator : RC/JU
 Sample : Q1248-07
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A6333

Quant Time: Feb 03 08:31:49 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 31 11:40:55 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	365599	20.000	ng/u1	0.02
20) Naphthalene-d8	10.557	136	1532078	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.398	164	992419	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.181	188	2024956	20.000	ng/u1	-0.03
79) Chrysene-d12	21.622	240	2054378	20.000	ng/u1	-0.02
88) Perylene-d12	24.992	264	2153028	20.000	ng/u1	-0.03

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	25183	2.862	ng/uL	0.01
4) Pyridine-d5	3.711	84	352319	13.876	ng/u1	0.02
7) Phenol-d5	6.963	99	499834	15.449	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	7.116	67	291618	17.028	ng/u1	0.02
11) 2-Chlorophenol-d4	7.328	132	408231	16.068	ng/u1	0.02
15) 4-Methylphenol-d8	8.487	113	418113	15.820	ng/u1	0.02
21) Nitrobenzene-d5	8.922	128	191660	15.636	ng/u1	0.01
24) 2-Nitrophenol-d4	9.640	143	171136	12.939	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.181	165	372087	15.335	ng/u1	0.00
31) 4-Chloroaniline-d4	10.687	131	538344	14.836	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	1309962	17.696	ng/u1	0.00
49) Acenaphthylene-d8	14.092	160	976548	11.697	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	183311	12.469	ng/u1	0.00
60) Fluorene-d10	15.398	176	615885	9.880	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.516	200	64904	5.195	ng/u1	-0.02
73) Anthracene-d10	17.286	188	844653	8.495	ng/u1	-0.03
81) Pyrene-d10	19.645	212	746940	6.786	ng/u1	-0.03
92) Benzo(a)pyrene-d12	24.745	264	478176	4.257	ng/u1	-0.04

Target Compounds	Qvalue

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

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