

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
 Data File : BP023640.D
 Acq On : 17 Jan 2025 23:25
 Operator : RC/JU
 Sample : Q1088-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B00

Quant Time: Jan 17 23:58:53 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP011425.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 14 16:27:27 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	418377	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1822538	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	1178013	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.239	188	2427791	20.000	ng/u1	0.00
79) Chrysene-d12	21.698	240	2511268	20.000	ng/u1	-0.01
88) Perylene-d12	25.116	264	2750805	20.000	ng/u1	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	57859	5.099	ng/uL	0.00
4) Pyridine-d5	3.717	84	264976	8.377	ng/u1	0.00
7) Phenol-d5	6.964	99	321548	8.927	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	734876	34.827	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	870750	30.611	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	598364	21.417	ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	481214	34.593	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	390810	42.801	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	876437	31.557	ng/u1	0.00
31) 4-Chloroaniline-d4	10.704	131	1166357	26.306	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	3269921	35.667	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	3711802	34.034	ng/u1	0.00
54) 4-Nitrophenol-d4	14.604	143	112961	7.413	ng/u1	0.00
60) Fluorene-d10	15.445	176	2958691	37.290	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.545	200	251053	31.781	ng/u1	0.00
73) Anthracene-d10	17.339	188	5201503	39.689	ng/u1	0.00
81) Pyrene-d10	19.710	212	6156446	44.252	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.886	264	6508723	43.539	ng/u1	-0.02

Target Compounds	Qvalue

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

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