

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
 Data File : BP023644.D
 Acq On : 18 Jan 2025 02:09
 Operator : RC/JU
 Sample : Q1088-07
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B02

Quant Time: Jan 18 02:56:01 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	411997	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	1811545	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	1179321	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2352308	20.000	ng/u1	0.00
79) Chrysene-d12	21.686	240	2335743	20.000	ng/u1	-0.02
88) Perylene-d12	25.098	264	2461900	20.000	ng/u1	-0.05
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	69389	6.210	ng/uL	0.00
4) Pyridine-d5	3.717	84	268516	8.620	ng/u1	0.00
7) Phenol-d5	6.963	99	409752	11.552	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	952052	45.819	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	1116297	39.851	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	784902	28.529	ng/u1	0.00
21) Nitrobenzene-d5	8.952	128	632089	45.714	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	557999	61.482	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	1210149	43.837	ng/u1	0.00
31) 4-Chloroaniline-d4	10.704	131	1322298	30.004	ng/u1	0.00
46) Dimethylphthalate-d6	13.834	166	4262422	46.441	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	4853501	44.452	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	166255	10.898	ng/u1	0.00
60) Fluorene-d10	15.439	176	3706294	46.660	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.545	200	342386	44.734	ng/u1	0.00
73) Anthracene-d10	17.339	188	6098319	48.026	ng/u1	0.00
81) Pyrene-d10	19.710	212	7030466	54.332	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.868	264	7226807	54.016	ng/u1	-0.04

Target Compounds Qvalue

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

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