

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011725\
 Data File : BP023646.D
 Acq On : 18 Jan 2025 03:31
 Operator : RC/JU
 Sample : Q1088-09
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AM2

Quant Time: Jan 18 03:54:46 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	7.804	152	563144	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	2401511	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	1516083	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	2968438	20.000	ng/u1	0.00
79) Chrysene-d12	21.686	240	2759255	20.000	ng/u1	-0.02
88) Perylene-d12	25.109	264	2952061	20.000	ng/u1	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	85437	5.594	ng/uL	0.00
4) Pyridine-d5	3.722	84	214126	5.029	ng/u1	0.00
7) Phenol-d5	6.963	99	491221	10.132	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	1282176	45.144	ng/u1	0.00
11) 2-Chlorophenol-d4	7.340	132	1518830	39.668	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	985023	26.193	ng/u1	0.00
21) Nitrobenzene-d5	8.946	128	864235	47.149	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	759397	63.117	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	1576951	43.091	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	605216	10.359	ng/u1	0.00
46) Dimethylphthalate-d6	13.839	166	5342187	45.277	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	6234279	44.416	ng/u1	0.00
54) 4-Nitrophenol-d4	14.604	143	151240	7.711	ng/u1	0.00
60) Fluorene-d10	15.439	176	4752896	46.545	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.545	200	423081	43.804	ng/u1	0.00
73) Anthracene-d10	17.333	188	7527422	46.976	ng/u1	-0.01
81) Pyrene-d10	19.710	212	8430827	55.154	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.874	264	8146227	50.778	ng/u1	-0.04

Target Compounds	Qvalue

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

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