

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012025\
 Data File : BP023656.D
 Acq On : 20 Jan 2025 15:54
 Operator : RC/JU
 Sample : Q1088-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B10

Quant Time: Jan 20 17:16:55 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	818499	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	3227635	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	1937739	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.240	188	3761145	20.000	ng/u1	0.00
79) Chrysene-d12	21.669	240	3899113	20.000	ng/u1	-0.04
88) Perylene-d12	25.074	264	4339464	20.000	ng/u1	-0.07
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	97578	4.396	ng/uL	0.00
4) Pyridine-d5	3.723	84	509539	8.234	ng/u1	0.00
7) Phenol-d5	6.964	99	612648	8.694	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	1369519	33.176	ng/u1	0.00
11) 2-Chlorophenol-d4	7.346	132	1765862	31.731	ng/u1	0.00
15) 4-Methylphenol-d8	8.493	113	1139208	20.843	ng/u1	0.00
21) Nitrobenzene-d5	8.952	128	925550	37.570	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	861897	53.301	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	1727930	35.131	ng/u1	0.00
31) 4-Chloroaniline-d4	10.705	131	2045646	26.052	ng/u1	0.00
46) Dimethylphthalate-d6	13.840	166	5356179	35.517	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	6230201	34.728	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	208521	8.319	ng/u1	0.00
60) Fluorene-d10	15.445	176	4749095	36.388	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	512254	41.859	ng/u1	0.00
73) Anthracene-d10	17.334	188	7631074	37.586	ng/u1	-0.01
81) Pyrene-d10	19.698	212	8838534	40.918	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.845	264	9700076	41.132	ng/u1	-0.06

Target Compounds	Qvalue

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

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