

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021025\
 Data File : BP024012.D
 Acq On : 10 Feb 2025 14:30
 Operator : RC/JU
 Sample : Q1336-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29C5

Quant Time: Feb 19 14:13:19 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 13 12:56:03 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	340584	20.000	ng/u1	0.00
20) Naphthalene-d8	10.534	136	1527390	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.381	164	967142	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	1870717	20.000	ng/u1	0.02
79) Chrysene-d12	21.610	240	1670985	20.000	ng/u1	0.00
88) Perylene-d12	24.957	264	1632864	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	44017	5.370	ng/uL	0.00
4) Pyridine-d5	3.705	84	284776	12.040	ng/u1	0.00
7) Phenol-d5	6.946	99	299956	9.952	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.099	67	589195	36.931	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	733132	30.975	ng/u1	0.00
15) 4-Methylphenol-d8	8.463	113	585521	23.781	ng/u1	0.00
21) Nitrobenzene-d5	8.905	128	416859	34.113	ng/u1	0.00
24) 2-Nitrophenol-d4	9.622	143	418586	31.745	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.163	165	777021	32.123	ng/u1	0.00
31) 4-Chloroaniline-d4	10.669	131	1164272	32.184	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2476856	34.334	ng/u1	0.00
49) Acenaphthylene-d8	14.075	160	2921672	35.909	ng/u1	0.00
54) 4-Nitrophenol-d4	14.610	143	64132	4.476	ng/u1	0.01
60) Fluorene-d10	15.387	176	2271296	37.388	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	224476	19.450	ng/u1	0.01
73) Anthracene-d10	17.287	188	3656231	39.806	ng/u1	0.02
81) Pyrene-d10	19.657	212	3978010	44.430	ng/u1	0.02
92) Benzo(a)pyrene-d12	24.721	264	3388558	39.774	ng/u1	0.00

Target Compounds Qvalue

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

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