

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021025\
 Data File : BP024021.D
 Acq On : 10 Feb 2025 21:14
 Operator : RC/JU
 Sample : Q1336-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29D2

Quant Time: Feb 12 15:04:17 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 03 11:19:36 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	387273	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.534	136	1665911	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.381	164	1019563	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.163	188	2037415	20.000	ng/u1	-0.02
79) Chrysene-d12	21.598	240	1944534	20.000	ng/u1	-0.04
88) Perylene-d12	24.945	264	1928884	20.000	ng/u1	-0.06
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	51126	5.485	ng/uL	0.00
4) Pyridine-d5	3.705	84	252575	9.391	ng/u1	0.00
7) Phenol-d5	6.946	99	275993	8.053	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.099	67	579018	31.918	ng/u1	0.00
11) 2-Chlorophenol-d4	7.310	132	755576	28.074	ng/u1	0.00
15) 4-Methylphenol-d8	8.469	113	521235	18.618	ng/u1	0.00
21) Nitrobenzene-d5	8.910	128	402154	30.173	ng/u1	0.00
24) 2-Nitrophenol-d4	9.622	143	411853	28.638	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.163	165	761633	28.868	ng/u1	0.00
31) 4-Chloroaniline-d4	10.669	131	1045903	26.508	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2388222	31.404	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	2752380	32.089	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.604	143	80627	5.338	ng/u1	0.00
60) Fluorene-d10	15.386	176	2206407	34.452	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	262365	20.873	ng/u1	0.00
73) Anthracene-d10	17.269	188	3608687	36.074	ng/u1	-0.02
81) Pyrene-d10	19.633	212	4101778	39.368	ng/u1	-0.03
92) Benzo(a)pyrene-d12	24.721	264	3773864	37.499	ng/u1	-0.05

Target Compounds Qvalue

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

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