

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021025\
 Data File : BP024010.D
 Acq On : 10 Feb 2025 13:09
 Operator : RC/JU
 Sample : Q1329-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHQJ3

Quant Time: Feb 12 14:30:42 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	377698	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.534	136	1683625	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.380	164	1061290	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.168	188	2014033	20.000	ng/u1	-0.02
79) Chrysene-d12	21.598	240	1757157	20.000	ng/u1	-0.04
88) Perylene-d12	24.956	264	1743059	20.000	ng/u1	-0.05
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	47948	5.275	ng/uL	0.00
4) Pyridine-d5	3.704	84	180792	6.892	ng/u1	0.00
7) Phenol-d5	6.951	99	274314	8.207	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.098	67	571621	32.309	ng/u1	0.00
11) 2-Chlorophenol-d4	7.304	132	741832	28.262	ng/u1	-0.01
15) 4-Methylphenol-d8	8.469	113	527974	19.336	ng/u1	0.00
21) Nitrobenzene-d5	8.904	128	410888	30.504	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.622	143	420424	28.926	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.163	165	777900	29.175	ng/u1	0.00
31) 4-Chloroaniline-d4	10.669	131	996819	24.998	ng/u1	0.00
46) Dimethylphthalate-d6	13.780	166	2594072	32.769	ng/u1	-0.01
49) Acenaphthylene-d8	14.069	160	2950972	33.052	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.604	143	62607	3.982	ng/u1	0.00
60) Fluorene-d10	15.386	176	2299325	34.491	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	214173	17.237	ng/u1	0.00
73) Anthracene-d10	17.274	188	3721785	37.636	ng/u1	-0.01
81) Pyrene-d10	19.639	212	3950094	41.955	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.727	264	3494750	38.427	ng/u1	-0.05

Target Compounds Qvalue

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

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