

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021025\
 Data File : BP024014.D
 Acq On : 10 Feb 2025 15:51
 Operator : RC/JU
 Sample : Q1336-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29C6

Quant Time: Feb 12 14:37:13 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	343102	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.540	136	1497923	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.381	164	930384	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.169	188	1872060	20.000	ng/u1	-0.02
79) Chrysene-d12	21.616	240	1645493	20.000	ng/u1	-0.02
88) Perylene-d12	24.963	264	1677341	20.000	ng/u1	-0.05

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	42059	5.093	ng/uL	0.00
4) Pyridine-d5	3.705	84	151617	6.363	ng/u1	0.00
7) Phenol-d5	6.952	99	278001	9.156	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.099	67	615059	38.269	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	777909	32.625	ng/u1	0.00
15) 4-Methylphenol-d8	8.469	113	547613	22.078	ng/u1	0.00
21) Nitrobenzene-d5	8.910	128	434483	36.254	ng/u1	0.00
24) 2-Nitrophenol-d4	9.628	143	442010	34.181	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	833291	35.126	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	1132763	31.929	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2633705	37.951	ng/u1	0.00
49) Acenaphthylene-d8	14.075	160	2973700	37.993	ng/u1	0.00
54) 4-Nitrophenol-d4	14.604	143	73314	5.320	ng/u1	0.00
60) Fluorene-d10	15.387	176	2384259	40.798	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.498	200	271080	23.472	ng/u1	-0.01
73) Anthracene-d10	17.275	188	3849700	41.882	ng/u1	-0.01
81) Pyrene-d10	19.651	212	4370073	49.565	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.745	264	3896070	44.518	ng/u1	-0.03

Target Compounds Qvalue

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

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