

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021225\
 Data File : BP024054.D
 Acq On : 12 Feb 2025 17:43
 Operator : RC/JU
 Sample : P5368-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YE626

Quant Time: Feb 14 12:25:35 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	7.757	152	562995	20.000	ng/u1	0.00
20) Naphthalene-d8	10.534	136	2156856	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.386	164	1187964	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.180	188	2080003	20.000	ng/u1	0.01
79) Chrysene-d12	21.621	240	2099140	20.000	ng/u1	0.02
88) Perylene-d12	24.974	264	2389249	20.000	ng/u1	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	213	0.016	ng/uL	0.01
4) Pyridine-d5	3.693	84	213	0.005	ng/u1	0.00
7) Phenol-d5	6.951	99	98	0.002	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.093	67	2978	0.113	ng/u1	0.00
11) 2-Chlorophenol-d4	7.281	132	85	0.002	ng/u1	-0.02
15) 4-Methylphenol-d8	8.469	113	249	0.006	ng/u1	0.00
21) Nitrobenzene-d5	8.928	128	1735	0.101	ng/u1	0.02
24) 2-Nitrophenol-d4	9.616	143	49	0.003	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	187	0.005	ng/u1	0.01
31) 4-Chloroaniline-d4	10.645	131	16504	0.323	ng/u1	-0.02
46) Dimethylphthalate-d6	13.786	166	267	0.003	ng/u1	0.00
49) Acenaphthylene-d8	14.080	160	785	0.008	ng/u1	0.01
54) 4-Nitrophenol-d4	14.604	143	2191	0.125	ng/u1	0.00
60) Fluorene-d10	15.392	176	268	0.004	ng/u1	0.01
65) 4,6-Dinitro-2-methylph...	15.392	200	1104	0.086	ng/u1	-0.11
73) Anthracene-d10	0.000	188	0d	0.000	ng/u1	
81) Pyrene-d10	19.663	212	2736	0.024	ng/u1	0.02
92) Benzo(a)pyrene-d12	24.745	264	2283	0.018	ng/u1	0.02

Target Compounds Qvalue

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

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