

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
 Data File : BP023981.D
 Acq On : 06 Feb 2025 22:05
 Operator : RC/JU
 Sample : Q1229-09
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29A9

Quant Time: Feb 07 07:34:20 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	7.769	152	367393	20.000	ng/u1	0.00
20) Naphthalene-d8	10.540	136	1577382	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.398	164	986278	20.000	ng/u1	0.01
64) Phenanthrene-d10	17.198	188	2008927	20.000	ng/u1	0.01
79) Chrysene-d12	21.639	240	2008209	20.000	ng/u1	0.00
88) Perylene-d12	25.015	264	2119424	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	54389	6.151	ng/uL	0.00
4) Pyridine-d5	3.705	84	258553	10.133	ng/u1	0.00
7) Phenol-d5	6.957	99	376213	11.571	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	755399	43.894	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	1023159	40.074	ng/u1	0.00
15) 4-Methylphenol-d8	8.475	113	694042	26.131	ng/u1	0.00
21) Nitrobenzene-d5	8.916	128	560509	44.414	ng/u1	0.00
24) 2-Nitrophenol-d4	9.628	143	604853	44.418	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.175	165	1060924	42.469	ng/u1	0.00
31) 4-Chloroaniline-d4	10.675	131	1331311	35.635	ng/u1	0.00
46) Dimethylphthalate-d6	13.804	166	3334801	45.330	ng/u1	0.01
49) Acenaphthylene-d8	14.087	160	3829940	46.159	ng/u1	0.00
54) 4-Nitrophenol-d4	14.622	143	144709	9.905	ng/u1	0.01
60) Fluorene-d10	15.416	176	2974147	48.007	ng/u1	0.02
65) 4,6-Dinitro-2-methylph...	15.528	200	447412	36.100	ng/u1	0.02
73) Anthracene-d10	17.292	188	5036117	51.056	ng/u1	0.00
81) Pyrene-d10	19.669	212	5847127	54.340	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.768	264	5912129	53.464	ng/u1	0.00

Target Compounds Qvalue

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

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