

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
 Data File : BP023988.D
 Acq On : 07 Feb 2025 03:29
 Operator : RC/JU
 Sample : Q1229-12
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29B8

Quant Time: Feb 07 07:35:18 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	432494	20.000	ng/u1	0.00
20) Naphthalene-d8	10.540	136	1897428	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.393	164	1214524	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.192	188	2378708	20.000	ng/u1	0.00
79) Chrysene-d12	21.627	240	2256720	20.000	ng/u1	-0.01
88) Perylene-d12	24.992	264	2376267	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	56070	5.387	ng/uL	0.00
4) Pyridine-d5	3.705	84	267318	8.900	ng/u1	0.00
7) Phenol-d5	6.952	99	475356	12.420	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	805300	39.750	ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	1101894	36.661	ng/u1	0.00
15) 4-Methylphenol-d8	8.469	113	912112	29.173	ng/u1	0.00
21) Nitrobenzene-d5	8.910	128	623965	41.103	ng/u1	0.00
24) 2-Nitrophenol-d4	9.628	143	695062	42.433	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.169	165	1214427	40.414	ng/u1	0.00
31) 4-Chloroaniline-d4	10.669	131	1469124	32.691	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	3986351	44.003	ng/u1	0.00
49) Acenaphthylene-d8	14.081	160	4345141	42.527	ng/u1	0.00
54) 4-Nitrophenol-d4	14.616	143	160960	8.947	ng/u1	0.00
60) Fluorene-d10	15.398	176	3523362	46.184	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.516	200	518661	35.344	ng/u1	0.00
73) Anthracene-d10	17.286	188	5634311	48.241	ng/u1	0.00
81) Pyrene-d10	19.651	212	6278143	51.920	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.768	264	6285746	50.699	ng/u1	0.00

Target Compounds	Qvalue

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

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