

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020725\
 Data File : BP024001.D
 Acq On : 07 Feb 2025 13:49
 Operator : RC/JU
 Sample : Q1229-16
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E29C2

Quant Time: Feb 12 12:52:25 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	403258	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.540	136	1723703	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.381	164	1039163	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	2070439	20.000	ng/u1	0.00
79) Chrysene-d12	21.616	240	1987996	20.000	ng/u1	-0.02
88) Perylene-d12	24.986	264	2067583	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.287	96	43084	4.439	ng/uL	0.00
4) Pyridine-d5	3.705	84	308074	11.000	ng/u1	0.00
7) Phenol-d5	6.952	99	282418	7.914	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.099	67	621323	32.892	ng/u1	0.00
11) 2-Chlorophenol-d4	7.305	132	795420	28.383	ng/u1	-0.01
15) 4-Methylphenol-d8	8.469	113	579643	19.883	ng/u1	0.00
21) Nitrobenzene-d5	8.910	128	449328	32.582	ng/u1	0.00
24) 2-Nitrophenol-d4	9.622	143	471163	31.663	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.163	165	878463	32.180	ng/u1	0.00
31) 4-Chloroaniline-d4	10.669	131	674717	16.527	ng/u1	0.00
46) Dimethylphthalate-d6	13.787	166	2747960	35.452	ng/u1	0.00
49) Acenaphthylene-d8	14.069	160	3181034	36.387	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.610	143	86223	5.601	ng/u1	0.00
60) Fluorene-d10	15.387	176	2466225	37.783	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.510	200	328561	25.723	ng/u1	0.00
73) Anthracene-d10	17.281	188	3989535	39.244	ng/u1	0.00
81) Pyrene-d10	19.657	212	4522929	42.461	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.739	264	4297366	39.836	ng/u1	-0.04

Target Compounds Qvalue

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

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