

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
 Data File : BP012565.D
 Acq On : 08 Nov 2022 23:15
 Operator : CG/JU
 Sample : N5407-12
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4Z1

Quant Time: Nov 09 00:38:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 08 22:40:23 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	244794	20.000	ng/u1	0.00
20) Naphthalene-d8	10.569	136	1245449	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.422	164	894900	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.186	188	2009541	20.000	ng/u1	0.00
79) Chrysene-d12	21.286	240	1821626	20.000	ng/u1	# 0.00
88) Perylene-d12	23.663	264	1650811	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	23664	3.892	ng/uL	0.00
4) Pyridine-d5	3.652	84	145205	8.314	ng/u1	0.00
7) Phenol-d5	6.963	99	181305	8.258	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	392415	29.939	ng/u1	0.00
11) 2-Chlorophenol-d4	7.305	132	415499	25.914	ng/u1	0.00
15) 4-Methylphenol-d8	8.499	113	331332	19.080	ng/u1	0.00
21) Nitrobenzene-d5	8.940	128	286697	35.498	ng/u1	0.00
24) 2-Nitrophenol-d4	9.657	143	304386	36.311	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	529697	29.157	ng/u1	0.00
31) 4-Chloroaniline-d4	10.722	131	541678	20.025	ng/u1	0.00
46) Dimethylphthalate-d6	13.839	166	2125733	34.700	ng/u1	0.00
49) Acenaphthylene-d8	14.116	160	2215972	31.653	ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	73117	6.306	ng/u1	0.01
60) Fluorene-d10	15.422	176	1841784	35.115	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	352395	33.682	ng/u1	0.00
73) Anthracene-d10	17.286	188	3183687	36.769	ng/u1	0.00
81) Pyrene-d10	19.527	212	3730841	40.793	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.510	264	3159858	39.175	ng/u1	0.00

Target Compounds Qvalue

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

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