

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112822\
 Data File : BP012832.D
 Acq On : 29 Nov 2022 00:29
 Operator : CG/JU
 Sample : N5666-13
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHB35

Quant Time: Nov 29 00:57:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Nov 27 22:55:08 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	628226	20.000	ng/u1	0.00
20) Naphthalene-d8	10.828	136	2825410	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1894589	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	3803336	20.000	ng/u1	0.00
79) Chrysene-d12	21.486	240	2294199	20.000	ng/u1	-0.01
88) Perylene-d12	23.998	264	1928612	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	64350	4.389	ng/uL	0.00
4) Pyridine-d5	3.822	84	381089	8.686	ng/u1	0.00
7) Phenol-d5	7.163	99	509969	9.879	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	533654	16.776	ng/u1	0.00
11) 2-Chlorophenol-d4	7.540	132	558148	14.189	ng/u1	0.00
15) 4-Methylphenol-d8	8.710	113	502869	12.090	ng/u1	-0.01
21) Nitrobenzene-d5	9.181	128	290238	16.783	ng/u1	0.00
24) 2-Nitrophenol-d4	9.904	143	291666	15.884	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	596378	13.920	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	739235	12.600	ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	2170790	16.646	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	2423279	15.926	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	106910	4.598	ng/u1	0.00
60) Fluorene-d10	15.639	176	1927443	16.877	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	99265	5.461	ng/u1	0.00
73) Anthracene-d10	17.504	188	2997823	17.906	ng/u1	0.00
81) Pyrene-d10	19.727	212	2918810	21.896	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.833	264	1639078	17.302	ng/u1	-0.01

Target Compounds Qvalue

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

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