

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012466.D
 Acq On : 03 Nov 2022 00:02
 Operator : CG/JU
 Sample : N5300-13
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA91

Quant Time: Nov 03 02:28:36 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 03 00:09:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	271186	20.000	ng/u1	0.00
20) Naphthalene-d8	10.587	136	1272303	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.434	164	876477	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	1935066	20.000	ng/u1	0.00
79) Chrysene-d12	21.292	240	1896600	20.000	ng/u1	0.00
88) Perylene-d12	23.674	264	1709656	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	23614	3.506	ng/uL	0.00
4) Pyridine-d5	3.664	84	102335	5.289	ng/u1	0.00
7) Phenol-d5	6.963	99	166190	6.833	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	426881	29.399	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	449705	25.318	ng/u1	0.00
15) 4-Methylphenol-d8	8.504	113	330019	17.155	ng/u1	0.00
21) Nitrobenzene-d5	8.952	128	311090	37.705	ng/u1	0.00
24) 2-Nitrophenol-d4	9.675	143	339792	39.679	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	617077	33.250	ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	749340	27.117	ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	2422118	40.369	ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	2460094	35.879	ng/u1	0.00
54) 4-Nitrophenol-d4	14.675	143	82768	7.289	ng/u1	0.01
60) Fluorene-d10	15.434	176	2070990	40.315	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	309374	30.708	ng/u1	0.00
73) Anthracene-d10	17.298	188	3523302	42.257	ng/u1	0.00
81) Pyrene-d10	19.539	212	4183953	43.939	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.527	264	3657587	43.784	ng/u1	0.00

Target Compounds Qvalue

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

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