

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012452.D
 Acq On : 02 Nov 2022 14:12
 Operator : CG/JU
 Sample : N5270-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA66

Quant Time: Nov 02 23:14:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	250744	20.000	ng/ul	0.00
20) Naphthalene-d8	10.587	136	1176344	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.439	164	789901	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1768285	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	1591955	20.000	ng/ul	0.00
88) Perylene-d12	23.680	264	1397326	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	27416	4.403	ng/uL	0.00
4) Pyridine-d5	3.664	84	111792	6.249	ng/ul	0.00
7) Phenol-d5	6.963	99	172116	7.653	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	488126	36.358	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	485219	29.544	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	350005	19.677	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	331046	43.397	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	358225	45.244	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	629197	36.668	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	666136	26.073	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	2263886	41.867	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	2503376	40.512	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	57578	5.626	ng/ul	0.00
60) Fluorene-d10	15.439	176	1973218	42.622	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	297039	32.264	ng/ul	0.00
73) Anthracene-d10	17.304	188	3219104	42.250	ng/ul	0.00
81) Pyrene-d10	19.545	212	3677413	46.010	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.527	264	2967410	43.462	ng/ul	-0.02

Target Compounds Qvalue

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

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