

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
 Data File : BP012478.D
 Acq On : 03 Nov 2022 07:27
 Operator : CG/JU
 Sample : N5321-23
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBWW5

Quant Time: Nov 03 07:55:15 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	283755	20.000 ng/u1	0.00
20) Naphthalene-d8	10.581	136	1337514	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.434	164	900684	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.198	188	1951399	20.000 ng/u1	0.00
79) Chrysene-d12	21.292	240	1842034	20.000 ng/u1	0.00
88) Perylene-d12	23.674	264	1639388	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.234	96	26175	3.714 ng/uL	0.00
4) Pyridine-d5	3.664	84	99624	4.921 ng/u1	0.00
7) Phenol-d5	6.958	99	213828	8.402 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	457878	30.137 ng/u1	0.00
11) 2-Chlorophenol-d4	7.311	132	511873	27.541 ng/u1	-0.01
15) 4-Methylphenol-d8	8.499	113	410725	20.405 ng/u1	0.00
21) Nitrobenzene-d5	8.952	128	320901	36.998 ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	365732	40.626 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	685854	35.154 ng/u1	0.00
31) 4-Chloroaniline-d4	10.734	131	482064	16.594 ng/u1	0.00
46) Dimethylphthalate-d6	13.851	166	2379377	38.591 ng/u1	0.00
49) Acenaphthylene-d8	14.128	160	2433153	34.532 ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	96352	8.257 ng/u1	0.00
60) Fluorene-d10	15.434	176	2029698	38.449 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	406911	40.051 ng/u1	-0.01
73) Anthracene-d10	17.298	188	3306436	39.324 ng/u1	0.00
81) Pyrene-d10	19.539	212	3751858	40.568 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.522	264	3260504	40.704 ng/u1	0.00

Target Compounds Qvalue

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

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