

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110222\
 Data File : BF130934.D
 Acq On : 02 Nov 2022 17:05
 Operator : CG\JU
 Sample : N5396-17
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-H

Quant Time: Nov 03 04:33:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 03 03:37:09 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.922	152	67667	20.000	ng	0.00
21) Naphthalene-d8	8.210	136	256646	20.000	ng	# 0.00
39) Acenaphthene-d10	9.969	164	142079	20.000	ng	0.00
64) Phenanthrene-d10	11.463	188	256449	20.000	ng	0.00
76) Chrysene-d12	14.110	240	143959	20.000	ng	-0.01
86) Perylene-d12	15.621	264	127595	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.551	112	474619	121.596	ng	0.01
7) Phenol-d6	6.545	99	538245	106.116	ng	0.00
23) Nitrobenzene-d5	7.486	82	463364	108.867	ng	0.00
42) 2,4,6-Tribromophenol	10.763	330	207203	173.244	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	843113	89.717	ng	0.00
79) Terphenyl-d14	13.057	244	843111	107.320	ng	0.00

Target Compounds Qvalue

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

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