

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111122\
 Data File : BN022606.D
 Acq On : 11 Nov 2022 10:21
 Operator : CG/JU
 Sample : PB148932BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB148932BL

Quant Time: Nov 11 22:14:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN111122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 23:28:29 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	1747	0.400	ng	0.00
7) Naphthalene-d8	10.731	136	4841	0.400	ng	0.01
13) Acenaphthene-d10	14.582	164	2628	0.400	ng	0.01
19) Phenanthrene-d10	17.341	188	5036	0.400	ng	0.02
29) Chrysene-d12	21.546	240	4043	0.400	ng	0.00
35) Perylene-d12	23.987	264	3439	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.450	112	2043	0.404	ng	0.00
5) Phenol-d6	7.097	99	2293	0.363	ng	0.00
8) Nitrobenzene-d5	9.097	82	1550	0.337	ng	0.02
11) 2-Methylnaphthalene-d10	12.330	152	2630	0.333	ng	0.01
14) 2,4,6-Tribromophenol	16.100	330	457	0.360	ng	0.03
15) 2-Fluorobiphenyl	13.203	172	3830	0.358	ng	0.00
27) Fluoranthene-d10	19.381	212	4886	0.352	ng	0.00
31) Terphenyl-d14	19.976	244	5254	0.419	ng	0.00

Target Compounds	Qvalue

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

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