

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110423\
 Data File : BN028467.D
 Acq On : 04 Nov 2023 10:34
 Operator : MA/JU
 Sample : PB156879BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB156879BL

Quant Time: Nov 06 04:43:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN110123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 02 06:26:50 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.717	152	9344	0.400	ng	0.00
7) Naphthalene-d8	10.509	136	28629	0.400	ng	# 0.00
13) Acenaphthene-d10	14.363	164	14526	0.400	ng	0.00
19) Phenanthrene-d10	17.108	188	35608	0.400	ng	#-0.01
29) Chrysene-d12	21.328	240	24990	0.400	ng	0.00
35) Perylene-d12	23.648	264	8827	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.298	112	11002	0.451	ng	0.00
5) Phenol-d6	6.886	99	12729	0.412	ng	0.00
8) Nitrobenzene-d5	8.864	82	7508	0.406	ng	-0.01
11) 2-Methylnaphthalene-d10	12.106	152	17140	0.419	ng	0.00
14) 2,4,6-Tribromophenol	15.855	330	2847	0.503	ng	0.00
15) 2-Fluorobiphenyl	13.006	172	3195	0.185	ng	0.00
27) Fluoranthene-d10	19.157	212	34557	0.393	ng	0.00
31) Terphenyl-d14	19.770	244	25953	0.493	ng	0.00

Target Compounds Qvalue

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

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