

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
 Data File : BN016689.D
 Acq On : 03 Oct 2021 03:23
 Operator : CG/JU
 Sample : PB139503BL
 Misc :
 ALS Vial : 57 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB139503BL

Quant Time: Oct 04 03:36:33 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 01 13:35:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.642	152	29158	0.40	ng	0.00
7) Naphthalene-d8	10.407	136	127438	0.40	ng	0.00
13) Acenaphthene-d10	14.269	164	65675	0.40	ng	0.00
19) Phenanthrene-d10	17.030	188	124355	0.40	ng	0.00
29) Chrysene-d12	21.238	240	104653	0.40	ng	0.00
35) Perylene-d12	23.484	264	107260	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.282	112	25937	0.35	ng	0.03
5) Phenol-d6	6.831	99	22797	0.28	ng	0.00
8) Nitrobenzene-d5	8.784	82	23718	0.29	ng	0.00
11) 2-Methylnaphthalene-d10	12.003	152	72007	0.38	ng	0.00
14) 2,4,6-Tribromophenol	15.766	330	6461	0.22	ng	0.00
15) 2-Fluorobiphenyl	12.898	172	91061	0.34	ng	0.00
27) Fluoranthene-d10	19.068	212	125049	0.38	ng	0.00
31) Terphenyl-d14	19.678	244	95721	0.42	ng	0.00

Target Compounds	Qvalue

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

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