

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101221\
 Data File : BN016897.D
 Acq On : 12 Oct 2021 11:15
 Operator : CG/JU
 Sample : PB139115BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB139115BL

Quant Time: Oct 12 11:55:24 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN101221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 12 02:31:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.689	152	22284	0.400	ng	0.00
7) Naphthalene-d8	10.445	136	98496	0.400	ng	#-0.01
13) Acenaphthene-d10	14.294	164	50209	0.400	ng	0.00
19) Phenanthrene-d10	17.042	188	91439	0.400	ng	0.00
29) Chrysene-d12	21.227	240	83250	0.400	ng	#-0.01
35) Perylene-d12	23.478	264	85574	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.328	112	20206	0.354	ng	0.03
5) Phenol-d6	6.877	99	16664	0.260	ng	0.00
8) Nitrobenzene-d5	8.822	82	20613	0.310	ng	-0.01
11) 2-Methylnaphthalene-d10	12.038	152	56906	0.399	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	3397	0.140	ng	0.00
15) 2-Fluorobiphenyl	12.924	172	72003	0.375	ng	0.00
27) Fluoranthene-d10	19.073	212	94211	0.393	ng	0.00
31) Terphenyl-d14	19.679	244	81622	0.431	ng	0.00

Target Compounds Qvalue

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

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