

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032222\
 Data File : BN019090.D
 Acq On : 23 Mar 2022 14:08
 Operator : CG/JU
 Sample : PB143429BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB143429BL

Quant Time: Mar 23 16:45:55 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 17 13:18:23 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.854	152	4182	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	9788	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	5440	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	9717	0.400	ng	0.00
29) Chrysene-d12	21.422	240	7835	0.400	ng	# 0.00
35) Perylene-d12	23.793	264	6829	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.406	112	3002	0.320	ng	0.00
5) Phenol-d6	7.024	99	2064	0.210	ng	0.01
8) Nitrobenzene-d5	9.025	82	1925	0.273	ng	0.01
11) 2-Methylnaphthalene-d10	12.257	152	4756	0.293	ng	0.00
14) 2,4,6-Tribromophenol	15.995	330	388	0.148	ng	0.01
15) 2-Fluorobiphenyl	13.127	172	6714	0.308	ng	0.00
27) Fluoranthene-d10	19.265	212	10710	0.371	ng	0.00
31) Terphenyl-d14	19.864	244	6412	0.333	ng	0.00

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

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