

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062124\
 Data File : BN032147.D
 Acq On : 22 Jun 2024 01:20
 Operator : MA/JU
 Sample : PB161510BL
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB161510BL

Quant Time: Jun 22 01:51:50 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN061024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 19 11:30:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.639	152	9117	0.400	ng	0.00
7) Naphthalene-d8	10.421	136	29232	0.400	ng	0.00
13) Acenaphthene-d10	14.285	164	18728	0.400	ng	0.00
19) Phenanthrene-d10	17.029	188	36080	0.400	ng	# 0.00
29) Chrysene-d12	21.229	240	22024	0.400	ng	# 0.00
35) Perylene-d12	23.446	264	21874	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.248	112	9014	0.348	ng	0.00
5) Phenol-d6	6.830	99	9877	0.291	ng	0.00
8) Nitrobenzene-d5	8.798	82	7391	0.304	ng	0.00
11) 2-Methylnaphthalene-d10	12.013	152	15841	0.340	ng	0.00
14) 2,4,6-Tribromophenol	15.787	330	1581	0.176	ng	0.00
15) 2-Fluorobiphenyl	12.906	172	25350	0.357	ng	0.00
27) Fluoranthene-d10	19.068	212	30988	0.316	ng	0.00
31) Terphenyl-d14	19.667	244	22413	0.416	ng	0.00

Target Compounds Qvalue

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

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