

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072723\
 Data File : BN026767.D
 Acq On : 27 Jul 2023 16:28
 Operator : MA/JU
 Sample : PB154434BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB154434BL

Quant Time: Jul 27 18:32:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN062723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 27 18:29:30 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.885	152	1396	0.400	ng	# 0.04
7) Naphthalene-d8	10.703	136	3463	0.400	ng	# 0.06
13) Acenaphthene-d10	14.469	164	2537	0.400	ng	0.01
19) Phenanthrene-d10	17.244	188	4973	0.400	ng	0.02
29) Chrysene-d12	21.444	240	5090	0.400	ng	0.02
35) Perylene-d12	23.820	264	4651	0.400	ng	# 0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.437	112	1303	0.385	ng	-0.02
5) Phenol-d6	7.178	99	1579m	0.366	ng	0.11
8) Nitrobenzene-d5	9.230	82	1118m	0.377	ng	0.16
11) 2-Methylnaphthalene-d10	12.298	152	1902m	0.376	ng	0.06
14) 2,4,6-Tribromophenol	16.040	330	330	0.358	ng	0.05
15) 2-Fluorobiphenyl	13.144	172	3076m	0.314	ng	0.04
27) Fluoranthene-d10	19.267	212	6232	0.591	ng	0.01
31) Terphenyl-d14	19.853	244	5102	0.392	ng	0.00

Target Compounds Qvalue

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

