

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031225\
 Data File : BN036592.D
 Acq On : 12 Mar 2025 09:50
 Operator : RC/JU
 Sample : PB167057BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB167057BL

Quant Time: Mar 12 10:41:15 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN031025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 10 16:06:28 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.724	152	2081	0.400	ng	0.00
7) Naphthalene-d8	10.530	136	4397	0.400	ng	0.02
13) Acenaphthene-d10	14.377	164	2365	0.400	ng	0.01
19) Phenanthrene-d10	17.124	188	4163	0.400	ng	# 0.01
29) Chrysene-d12	21.304	240	2774	0.400	ng	# 0.00
35) Perylene-d12	23.560	264	2411	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.319	112	2126	0.438	ng	0.00
5) Phenol-d6	6.901	99	2057	0.343	ng	0.00
8) Nitrobenzene-d5	8.886	82	1694	0.354	ng	0.01
11) 2-Methylnaphthalene-d10	12.131	152	2268	0.347	ng	0.02
14) 2,4,6-Tribromophenol	15.882	330	200	0.186	ng	0.02
15) 2-Fluorobiphenyl	13.008	172	3233	0.235	ng	0.02
27) Fluoranthene-d10	19.150	212	4680	0.439	ng	0.00
31) Terphenyl-d14	19.750	244	2933	0.441	ng	0.00

Target Compounds Qvalue

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

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