

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040825\
 Data File : BN036860.D
 Acq On : 08 Apr 2025 13:45
 Operator : RC/JU
 Sample : Q1502-08DL 20X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Apr 08 14:18:29 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN031025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 04 17:30:43 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	2403	0.400	ng	0.00
7) Naphthalene-d8	10.466	136	6100	0.400	ng	#-0.01
13) Acenaphthene-d10	14.323	164	3298	0.400	ng	-0.01
19) Phenanthrene-d10	17.074	188	6675	0.400	ng	0.00
29) Chrysene-d12	21.268	240	5848	0.400	ng	0.00
35) Perylene-d12	23.510	264	6053	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.269	112	52744	9.418	ng	-0.01
5) Phenol-d6	6.850	99	67535	9.763	ng	-0.01
8) Nitrobenzene-d5	8.822	82	36163	5.450	ng	-0.02
11) 2-Methylnaphthalene-d10	0.000	152	0d	0.000	ng	
14) 2,4,6-Tribromophenol	15.820	330	13829	9.241	ng	-0.01
15) 2-Fluorobiphenyl	12.942	172	101987	5.316	ng	-0.02
27) Fluoranthene-d10	19.113	212	16	0.001	ng	0.00
31) Terphenyl-d14	19.712	244	83257	5.942	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.457	198	9388	5.007	ng	# 50
24) Pentachlorophenol	16.726	266	12444	5.404	ng	99

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

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