

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031425\
 Data File : BN036607.D
 Acq On : 14 Mar 2025 14:37
 Operator : RC/JU
 Sample : Q1502-08
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Mar 14 15:02:13 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	7.717	152	2909	0.400	ng	0.00
7) Naphthalene-d8	10.509	136	7981	0.400	ng	# 0.00
13) Acenaphthene-d10	14.366	164	4375	0.400	ng	0.00
19) Phenanthrene-d10	17.099	188	8959	0.400	ng	#-0.01
29) Chrysene-d12	21.304	240	5365	0.400	ng	0.00
35) Perylene-d12	23.554	264	3981	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.312	112	1233375	181.930	ng	0.00
5) Phenol-d6	6.901	99	1662926	198.573	ng	0.00
8) Nitrobenzene-d5	8.875	82	953549	109.829	ng	0.00
11) 2-Methylnaphthalene-d10	0.000	152	0d	0.000	ng	
14) 2,4,6-Tribromophenol	15.858	330	465620	234.541	ng	0.00
15) 2-Fluorobiphenyl	12.983	172	2628068	103.260	ng	0.00
27) Fluoranthene-d10	19.141	212	127	0.006	ng	0.00
31) Terphenyl-d14	19.745	244	1940216	150.943	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.499	198	435659	65.862	ng	# 21
24) Pentachlorophenol	16.764	266	457468	148.019	ng	99

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

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