

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091924\
 Data File : BN034042.D
 Acq On : 18 Sep 2024 19:13
 Operator : JU/RC
 Sample : P3845-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Sep 19 02:17:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 18 16:09:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.438	152	7056	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	21248	0.400	ng	# 0.00
13) Acenaphthene-d10	14.079	164	11207	0.400	ng	0.00
19) Phenanthrene-d10	16.840	188	24877	0.400	ng	# 0.00
29) Chrysene-d12	21.060	240	16205	0.400	ng	# 0.00
35) Perylene-d12	23.189	264	15627	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.076	112	2971330	148.381	ng	0.00
5) Phenol-d6	6.643	99	4028158	172.942	ng	0.00
8) Nitrobenzene-d5	8.579	82	2011619	123.766	ng	0.00
11) 2-Methylnaphthalene-d10	11.788	152	3	0.000	ng	0.00
14) 2,4,6-Tribromophenol	15.586	330	1184563	233.307	ng	0.00
15) 2-Fluorobiphenyl	12.700	172	4483178	98.356	ng	0.00
27) Fluoranthene-d10	18.906	212	290	0.005	ng	0.02
31) Terphenyl-d14	19.510	244	3766888	116.387	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.234	198	1241571	269.150	ng	# 51
24) Pentachlorophenol	16.493	266	494774	95.427	ng	100

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

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