

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091924\
 Data File : BN034043.D
 Acq On : 18 Sep 2024 19:49
 Operator : JU/RC
 Sample : P3845-08DL 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Sep 19 02:17:51 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	7922	0.400	ng	0.00
7) Naphthalene-d8	10.191	136	22980	0.400	ng	# 0.00
13) Acenaphthene-d10	14.079	164	11429	0.400	ng	0.00
19) Phenanthrene-d10	16.840	188	23143	0.400	ng	0.00
29) Chrysene-d12	21.060	240	13266	0.400	ng	# 0.00
35) Perylene-d12	23.189	264	12271	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.076	112	722073	32.117	ng	0.00
5) Phenol-d6	6.636	99	942308	36.034	ng	0.00
8) Nitrobenzene-d5	8.568	82	459778	26.156	ng	0.00
11) 2-Methylnaphthalene-d10	11.750	152	4	0.000	ng	-0.04
14) 2,4,6-Tribromophenol	15.587	330	224172	43.294	ng	0.00
15) 2-Fluorobiphenyl	12.700	172	1093393	23.522	ng	0.00
27) Fluoranthene-d10	18.915	212	58	0.001	ng	0.03
31) Terphenyl-d14	19.505	244	805433	30.399	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.234	198	186524	43.580	ng	# 23
24) Pentachlorophenol	16.493	266	74771	15.502	ng	100

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

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