

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041524\
 Data File : BN030945.D
 Acq On : 15 Apr 2024 18:58
 Operator : MA/JU
 Sample : P1908-08DL 50X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Apr 15 19:26:50 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN041524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 15 17:03:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.653	152	3595	0.400	ng	0.00
7) Naphthalene-d8	10.432	136	10824	0.400	ng	0.00
13) Acenaphthene-d10	14.295	164	6073	0.400	ng	0.00
19) Phenanthrene-d10	17.040	188	14444	0.400	ng	0.00
29) Chrysene-d12	21.240	240	11600	0.400	ng	0.00
35) Perylene-d12	23.457	264	11361	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.241	112	26455	3.303	ng	0.00
5) Phenol-d6	6.823	99	32945	3.396	ng	0.00
8) Nitrobenzene-d5	8.798	82	16267	2.171	ng	0.00
11) 2-Methylnaphthalene-d10	12.032	152	4	0.000	ng	0.00
14) 2,4,6-Tribromophenol	15.787	330	7948	3.373	ng	0.00
15) 2-Fluorobiphenyl	12.905	172	43927	2.152	ng	0.00
27) Fluoranthene-d10	19.084	212	4	0.000	ng	0.00
31) Terphenyl-d14	19.683	244	66455	2.468	ng	0.00
Target Compounds						Qvalue
20) 4,6-Dinitro-2-methylph...	15.428	198	4973	2.158	ng	# 54
24) Pentachlorophenol	16.693	266	9255	2.282	ng	99

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

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