

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040821\
 Data File : BN014239.D
 Acq On : 08 Apr 2021 14:24
 Operator : CG/JU
 Sample : M1557-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Apr 08 14:53:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN040821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 08 01:18:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.724	152	7653	0.40	ng	0.00
7) Naphthalene-d8	10.505	136	31105	0.40	ng	# 0.00
13) Acenaphthene-d10	14.346	164	18579	0.40	ng	-0.01
19) Phenanthrene-d10	17.092	188	38677	0.40	ng	#-0.01
29) Chrysene-d12	21.292	240	36635	0.40	ng	# 0.00
36) Perylene-d12	23.534	264	37206	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.324	112	2925473	143.37	ng	0.00
5) Phenol-d6	6.895	99	3816870	136.56	ng	0.00
8) Nitrobenzene-d5	8.870	82	2397356	111.46	ng	0.00
11) 2-Methylnaphthalene-d10	0.000	152	0d	0.00	ng	
14) 2,4,6-Tribromophenol	15.843	330	1828663	190.46	ng	0.00
15) 2-Fluorobiphenyl	12.976	172	5971006	92.40	ng	0.00
27) Fluoranthene-d10	0.000	212	0d	0.00	ng	
31) Terphenyl-d14	19.744	244	7247669	88.18	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.488	198	1707017	333.20	ng	Qvalue # 44
24) Pentachlorophenol	16.751	266	2015762	209.81	ng	96

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

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