

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101323\
 Data File : BN028276.D
 Acq On : 13 Oct 2023 12:01
 Operator : MA/JU
 Sample : 04346-08DL 50X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Oct 13 12:38:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN101223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 13 02:17:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.885	152	2838	0.400	ng	0.00
7) Naphthalene-d8	10.703	136	9324	0.400	ng	-0.01
13) Acenaphthene-d10	14.534	164	5775	0.400	ng	-0.01
19) Phenanthrene-d10	17.294	188	13200	0.400	ng	#-0.01
29) Chrysene-d12	21.497	240	12927	0.400	ng	0.00
35) Perylene-d12	23.955	264	14605	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.415	112	24436	2.773	ng	0.00
5) Phenol-d6	7.040	99	30891	2.847	ng	-0.01
8) Nitrobenzene-d5	9.080	82	15496	1.874	ng	-0.02
11) 2-Methylnaphthalene-d10	12.268	152	3	0.000	ng	-0.03
14) 2,4,6-Tribromophenol	16.040	330	8315	2.952	ng	-0.01
15) 2-Fluorobiphenyl	13.154	172	41882	2.061	ng	-0.01
27) Fluoranthene-d10	19.342	212	25	0.001	ng	0.00
31) Terphenyl-d14	19.918	244	65252	2.378	ng	0.00
Target Compounds						Qvalue
20) 4,6-Dinitro-2-methylph...	15.717	198	3802	2.577	ng	# 39
24) Pentachlorophenol	16.946	266	9752	3.052	ng	95

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

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