

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101321\
 Data File : BN016931.D
 Acq On : 13 Oct 2021 15:14
 Operator : CG/JU
 Sample : M3702-08DL 125X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Oct 13 16:12:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN101321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 13 02:19:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.679	152	15846	0.40	ng	0.00
7) Naphthalene-d8	10.445	136	69488	0.40	ng	#-0.01
13) Acenaphthene-d10	14.294	164	34624	0.40	ng	0.00
19) Phenanthrene-d10	17.030	188	67497	0.40	ng	#-0.01
29) Chrysene-d12	21.238	240	61169	0.40	ng	# 0.01
35) Perylene-d12	23.481	264	61668	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.282	112	58582	1.63	ng	0.00
5) Phenol-d6	6.859	99	61746	1.54	ng	0.00
8) Nitrobenzene-d5	8.822	82	42881	0.97	ng	0.00
11) 2-Methylnaphthalene-d10	12.034	152	6	0.00	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	22792	1.45	ng	0.00
15) 2-Fluorobiphenyl	12.924	172	137289	1.05	ng	0.00
27) Fluoranthene-d10	19.073	212	28	0.00	ng	0.00
31) Terphenyl-d14	19.679	244	164331	1.22	ng	0.00
Target Compounds						Qvalue
20) 4,6-Dinitro-2-methylph...	15.436	198	2395	0.36	ng	96
24) Pentachlorophenol	16.689	266	9277	0.61	ng	99

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

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