

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032321\
 Data File : BN014058.D
 Acq On : 24 Mar 2021 02:17
 Operator : CG/JU
 Sample : M1735-14
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D-20210316

Quant Time: Mar 24 02:59:42 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN032021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 23 11:06:27 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.700	152	5124	0.40	ng	0.00
7) Naphthalene-d8	10.467	136	19982	0.40	ng	#-0.01
13) Acenaphthene-d10	14.321	164	11792	0.40	ng	0.00
19) Phenanthrene-d10	17.055	188	26247	0.40	ng	-0.01
29) Chrysene-d12	21.239	240	25870	0.40	ng	# 0.00
36) Perylene-d12	23.441	264	23845	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.292	112	1330	0.11	ng	0.00
5) Phenol-d6	6.863	99	949	0.06	ng	0.00
8) Nitrobenzene-d5	8.845	82	3729	0.28	ng	0.00
11) 2-Methylnaphthalene-d10	12.062	152	9308	0.30	ng	0.00
14) 2,4,6-Tribromophenol	15.818	330	1098	0.29	ng	0.00
15) 2-Fluorobiphenyl	12.938	172	12783	0.29	ng	-0.01
27) Fluoranthene-d10	19.089	212	28224	0.39	ng	0.00
31) Terphenyl-d14	19.690	244	24099	0.42	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.239	88	27749	4.11	ng	96
6) bis(2-Chloroethyl)ether	7.129	93	881	0.07	ng	93
24) Pentachlorophenol	16.715	266	102	0.03	ng	97
34) Bis(2-ethylhexyl)phtha...	21.154	149	14668	0.53	ng	99

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

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