

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032321\
 Data File : BN014043.D
 Acq On : 23 Mar 2021 15:29
 Operator : CG/JU
 Sample : M1735-16
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 TT101D2-20210316

Quant Time: Mar 23 16:00:19 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	5106	0.40	ng	0.00
7) Naphthalene-d8	10.479	136	19570	0.40	ng	# 0.00
13) Acenaphthene-d10	14.320	164	11441	0.40	ng	0.00
19) Phenanthrene-d10	17.067	188	25389	0.40	ng	# 0.00
29) Chrysene-d12	21.239	240	25350	0.40	ng	# 0.00
36) Perylene-d12	23.448	264	23690	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.300	112	1718	0.15	ng	0.00
5) Phenol-d6	6.870	99	1252	0.08	ng	0.00
8) Nitrobenzene-d5	8.844	82	4247	0.33	ng	0.00
11) 2-Methylnaphthalene-d10	12.066	152	9953	0.33	ng	0.00
14) 2,4,6-Tribromophenol	15.817	330	1047	0.28	ng	0.00
15) 2-Fluorobiphenyl	12.950	172	13558	0.32	ng	0.00
27) Fluoranthene-d10	19.089	212	29454	0.42	ng	0.00
31) Terphenyl-d14	19.689	244	25579	0.46	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.238	88	13410	1.99	ng	# 88
6) bis(2-Chloroethyl)ether	7.128	93	459	0.03	ng	98
25) Phenanthrene	17.104	178	1447	0.02	ng	# 94
34) Bis(2-ethylhexyl)phtha...	21.154	149	4664	0.17	ng	98

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

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