

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032421\
 Data File : BN014075.D
 Acq On : 24 Mar 2021 17:59
 Operator : CG/JU
 Sample : M1787-06
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE-103D2-20210318

Quant Time: Mar 25 05:37:40 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	4980	0.40	ng	0.00
7) Naphthalene-d8	10.467	136	19172	0.40	ng	#-0.01
13) Acenaphthene-d10	14.321	164	11066	0.40	ng	0.00
19) Phenanthrene-d10	17.055	188	24227	0.40	ng	-0.01
29) Chrysene-d12	21.239	240	23034	0.40	ng	# 0.00
36) Perylene-d12	23.441	264	21750	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.292	112	1522	0.13	ng	0.00
5) Phenol-d6	6.863	99	1111	0.08	ng	0.00
8) Nitrobenzene-d5	8.832	82	3850	0.30	ng	-0.01
11) 2-Methylnaphthalene-d10	12.062	152	9766	0.33	ng	0.00
14) 2,4,6-Tribromophenol	15.805	330	884	0.24	ng	-0.01
15) 2-Fluorobiphenyl	12.938	172	13039	0.32	ng	-0.01
27) Fluoranthene-d10	19.089	212	27531	0.41	ng	0.00
31) Terphenyl-d14	19.685	244	25228	0.50	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.239	88	9460	1.44	ng	# 90
6) bis(2-Chloroethyl)ether	7.129	93	921	0.07	ng	# 61
25) Phenanthrene	17.104	178	1508	0.02	ng	# 95
34) Bis(2-ethylhexyl)phtha...	21.154	149	2166	0.09	ng	# 99

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

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