

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031021\
 Data File : BN013900.D
 Acq On : 10 Mar 2021 10:57
 Operator : CG/JU
 Sample : M1611-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWBR1-175-200-030921

Quant Time: Mar 10 11:28:15 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 08 10:32:09 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.717	152	5903	0.40	ng	0.00
7) Naphthalene-d8	10.492	136	23257	0.40	ng	#-0.01
13) Acenaphthene-d10	14.346	164	13078	0.40	ng	0.00
19) Phenanthrene-d10	17.092	188	28762	0.40	ng	# 0.00
29) Chrysene-d12	21.258	240	31685	0.40	ng	-0.01
36) Perylene-d12	23.487	264	29746	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.317	112	910	0.06	ng	0.00
5) Phenol-d6	6.887	99	691	0.03	ng	0.00
8) Nitrobenzene-d5	8.857	82	4701	0.31	ng	0.00
11) 2-Methylnaphthalene-d10	12.089	152	10434	0.28	ng	0.00
14) 2,4,6-Tribromophenol	15.831	330	1373	0.26	ng	-0.01
15) 2-Fluorobiphenyl	12.963	172	15475	0.33	ng	-0.01
27) Fluoranthene-d10	19.116	212	28471	0.34	ng	0.00
31) Terphenyl-d14	19.717	244	28400	0.41	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.255	88	18933	2.65	ng	95
9) Naphthalene	10.543	128	3561	0.06	ng	# 93
25) Phenanthrene	17.128	178	9444	0.13	ng	99
28) Fluoranthene	19.146	202	12229	0.14	ng	# 89
30) Pyrene	19.508	202	8715	0.09	ng	# 96
32) Benzo(a)anthracene	21.247	228	2199	0.02	ng	# 75
33) Chrysene	21.300	228	4169	0.04	ng	# 87
34) Bis(2-ethylhexyl)phtha...	21.184	149	18239	0.39	ng	97
37) Benzo(b)fluoranthene	22.819	252	4878	0.05	ng	# 1
39) Benzo(a)pyrene	23.387	252	2064	0.02	ng	# 1
41) Benzo(g,h,i)perylene	26.399	276	1924	0.02	ng	# 49

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

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