

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122220\
 Data File : BN013268.D
 Acq On : 22 Dec 2020 12:11
 Operator : CG/JU
 Sample : L5124-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 RE139D1-20201214

Quant Time: Dec 22 13:13:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN121620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 22 10:46:43 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.338	152	747	0.40	ng	0.00
7) Naphthalene-d8	10.112	136	2393	0.40	ng	# 0.01
13) Acenaphthene-d10	13.991	164	1427	0.40	ng	0.00
19) Phenanthrene-d10	16.763	188	3366	0.40	ng	# 0.00
29) Chrysene-d12	20.995	240	3776	0.40	ng	# 0.00
36) Perylene-d12	23.082	264	4325	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.962	112	313	0.10	ng	0.00
5) Phenol-d6	6.556	99	153	0.05	ng	0.02
8) Nitrobenzene-d5	8.540	82	601	0.24	ng	0.04
11) 2-Methylnaphthalene-d10	11.719	152	1009	0.24	ng	0.01
14) 2,4,6-Tribromophenol	15.526	330	231	0.26	ng	0.00
15) 2-Fluorobiphenyl	12.620	172	1547	0.26	ng	0.01
27) Fluoranthene-d10	18.811	212	4049	0.37	ng	0.00
31) Terphenyl-d14	19.426	244	3557	0.37	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.924	88	1529	1.17	ng	# 91
32) Benzo(a)anthracene	20.984	228	399	0.03	ng	# 68
33) Chrysene	21.037	228	449	0.03	ng	# 73
37) Benzo(b)fluoranthene	22.476	252	505	0.03	ng	# 1
38) Benzo(k)fluoranthene	22.517	252	471	0.03	ng	# 1
39) Benzo(a)pyrene	22.999	252	456	0.03	ng	# 1
40) Dibenzo(a,h)anthracene	25.149	278	342	0.02	ng	# 1
41) Benzo(g,h,i)perylene	25.758	276	510	0.03	ng	# 33

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

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