

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122220\
 Data File : BN013272.D
 Acq On : 22 Dec 2020 14:36
 Operator : CG/JU
 Sample : L5124-05
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DUP07-20201214

Quant Time: Dec 22 16:02:15 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	644	0.40	ng	0.00
7) Naphthalene-d8	10.112	136	1785	0.40	ng	# 0.01
13) Acenaphthene-d10	13.991	164	1207	0.40	ng	0.00
19) Phenanthrene-d10	16.775	188	2691	0.40	ng	# 0.01
29) Chrysene-d12	20.995	240	3076	0.40	ng	# 0.00
36) Perylene-d12	23.085	264	3476	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.962	112	316	0.12	ng	0.00
5) Phenol-d6	6.565	99	141	0.05	ng	0.02
8) Nitrobenzene-d5	8.540	82	630	0.34	ng	0.04
11) 2-Methylnaphthalene-d10	11.719	152	933	0.30	ng	0.01
14) 2,4,6-Tribromophenol	15.539	330	227	0.30	ng	0.01
15) 2-Fluorobiphenyl	12.608	172	1503	0.30	ng	0.00
27) Fluoranthene-d10	18.816	212	3582	0.40	ng	0.00
31) Terphenyl-d14	19.432	244	3081	0.39	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.925	88	707	0.63	ng	# 92
32) Benzo(a)anthracene	20.984	228	328	0.03	ng	# 64
33) Chrysene	21.038	228	385	0.03	ng	# 73
35) Indeno(1,2,3-cd)pyrene	25.153	276	392	0.02	ng	# 54
37) Benzo(b)fluoranthene	22.476	252	435	0.04	ng	# 1
38) Benzo(k)fluoranthene	22.517	252	411	0.03	ng	# 1
39) Benzo(a)pyrene	23.006	252	380	0.03	ng	# 1
40) Dibenzo(a,h)anthracene	25.156	278	308	0.02	ng	# 1
41) Benzo(g,h,i)perylene	25.755	276	562	0.04	ng	# 38

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

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