

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122220\
 Data File : BN013274.D
 Acq On : 22 Dec 2020 15:48
 Operator : CG/JU
 Sample : L5124-07
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 FB-20201214

Quant Time: Dec 22 17:03:19 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	626	0.40	ng	0.00
7) Naphthalene-d8	10.112	136	2170	0.40	ng	# 0.01
13) Acenaphthene-d10	13.991	164	1401	0.40	ng	0.00
19) Phenanthrene-d10	16.775	188	3142	0.40	ng	# 0.01
29) Chrysene-d12	21.003	240	3205	0.40	ng	# 0.00
36) Perylene-d12	23.086	264	3417	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	4.962	112	201	0.08	ng	0.00
5) Phenol-d6	6.581	99	118	0.05	ng	0.04
8) Nitrobenzene-d5	8.541	82	573	0.26	ng	0.04
11) 2-Methylnaphthalene-d10	11.711	152	1047	0.27	ng	0.00
14) 2,4,6-Tribromophenol	15.539	330	91	0.10	ng	0.01
15) 2-Fluorobiphenyl	12.608	172	1602	0.27	ng	0.00
27) Fluoranthene-d10	18.818	212	3788	0.37	ng	0.00
31) Terphenyl-d14	19.429	244	3159	0.39	ng	0.00
Target Compounds						Qvalue
32) Benzo(a)anthracene	20.992	228	312	0.03	ng	# 61
33) Chrysene	21.035	228	352	0.03	ng	# 71
35) Indeno(1,2,3-cd)pyrene	25.143	276	388	0.02	ng	# 54
37) Benzo(b)fluoranthene	22.477	252	426	0.04	ng	# 1
38) Benzo(k)fluoranthene	22.521	252	456	0.03	ng	# 1
40) Dibenzo(a,h)anthracene	25.160	278	285	0.02	ng	# 1
41) Benzo(g,h,i)perylene	25.763	276	571	0.04	ng	# 39

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

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