

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM081519\
 Data File : BM022151.D
 Acq On : 15 Aug 2019 19:46
 Operator : HP/JU
 Sample : K4242-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AF9

Quant Time: Aug 16 14:12:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 14 13:53:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	680	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.37	136	2924	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.25	164	1670	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.11	188	227382	0.40	ng/ul	0.07
16) Chrysene-d12	0.00	240	0	0.00	ng/ul	-21.23
20) Perylene-d12	23.42	264	4318	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	1363	0.28	ng/ul	-0.03
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds					Qvalue	
3) Naphthalene	10.42	128	425	0.051	ng/ul#	68
5) 2-Methylnaphthalene	12.06	142	395	0.072	ng/ul	99
7) Acenaphthylene	13.97	152	708	0.091	ng/ul#	72
8) Acenaphthene	14.31	153	374	0.059	ng/ul#	90
9) Fluorene	15.31	166	523	0.074	ng/ul#	96
21) Benzo(b)fluoranthene	22.76	252	72963	5.011	ng/ul	85
22) Benzo(k)fluoranthene	22.76	252	73148	4.439	ng/ul#	85
23) Benzo(a)pyrene	23.32	252	34928	2.359	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.62	276	29221	1.653	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.62	278	6594	0.461	ng/ul#	89
26) Benzo(g,h,i)perylene	26.30	276	23835	1.497	ng/ul	99

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

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