

Data Path : Z:\HPCHEM1\BNA_M\Data\BM122416\
 Data File : BM008626.D
 Acq On : 24 Dec 2016 10:31
 Operator : UM/SJ
 Sample : H6040-15
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA8W8

Quant Time: Dec 25 03:25:33 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 24 04:56:03 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	541	0.40	ng/ul	-0.09
2) Naphthalene-d8	10.43	136	1741	0.40	ng/ul	-0.07
6) Acenaphthene-d10	14.28	164	1056	0.40	ng/ul	-0.03
10) Phenanthrene-d10	17.14	188	71050	0.40	ng/ul	0.05
16) Chrysene-d12	21.41	240	59	0.40	ng/ul	0.14
20) Perylene-d12	23.72	264	108	0.40	ng/ul	0.22

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	409	0.15	ng/ul	-0.06
14) Fluoranthene-d10	19.17	212	131	0.00	ng/ul	0.06

Target Compounds				Qvalue		
3) Naphthalene	10.46	128	120	0.02	ng/ul#	2
5) 2-Methylnaphthalene	12.11	142	201	0.06	ng/ul	96
8) Acenaphthene	14.35	153	80	0.02	ng/ul#	74
17) Pyrene	19.63	202	457	2.00	ng/ul#	1
18) Benzo(a)anthracene	21.35	228	141	0.73	ng/ul#	1
19) Chrysene	21.47	228	219	1.03	ng/ul#	1
21) Benzo(b)fluoranthene	23.09	252	600	1.41	ng/ul#	1
22) Benzo(k)fluoranthene	23.09	252	702	1.69	ng/ul#	1
23) Benzo(a)pyrene	23.67	252	90	0.23	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.05	276	205	0.44	ng/ul#	37
25) Dibenzo(a,h)anthracene	26.05	278	342	0.91	ng/ul#	1
26) Benzo(g,h,i)perylene	26.61	276	106	0.26	ng/ul#	1

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

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