

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM112116\
 Data File : BM007962.D
 Acq On : 21 Nov 2016 17:32
 Operator : UM/SJ
 Sample : H5699-15
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4WE0

Quant Time: Nov 21 18:11:04 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 21 15:11:37 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	578	0.40	ng/ul	0.00
2) Naphthalene-d8	10.50	136	1827	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.35	164	1122	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.19	188	168785	0.40	ng/ul	0.09
16) Chrysene-d12	21.31	240	2389	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	135500	0.40	ng/ul	-0.16

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.12	152	727	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.14	212	2107	0.00	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.55	128	18	0.00	ng/ul#	1
5) 2-Methylnaphthalene	12.19	142	11	0.00	ng/ul#	71
7) Acenaphthylene	14.04	152	18	0.00	ng/ul#	1
8) Acenaphthene	14.42	153	5	0.00	ng/ul#	68
9) Fluorene	15.36	166	327	0.07	ng/ul#	30
11) Pentachlorophenol	17.00	266	1	0.00	ng/ul#	20
12) Phenanthrene	17.19	178	115	0.00	ng/ul#	1
13) Anthracene	17.38	178	7	0.00	ng/ul#	1
15) Fluoranthene	19.28	202	2	0.00	ng/ul#	1
17) Pyrene	19.50	202	367	0.04	ng/ul#	1
18) Benzo(a)anthracene	21.35	228	269	0.04	ng/ul#	52
19) Chrysene	21.35	228	278	0.03	ng/ul#	54
21) Benzo(b)fluoranthene	22.69	252	11	0.00	ng/ul#	1
22) Benzo(k)fluoranthene	22.79	252	13	0.00	ng/ul#	1
23) Benzo(a)pyrene	23.39	252	809	0.00	ng/ul#	45
24) Indeno(1,2,3-cd)pyrene	25.58	276	878	0.00	ng/ul#	51
25) Dibenzo(a,h)anthracene	25.65	278	1069	0.00	ng/ul#	1
26) Benzo(g,h,i)perylene	26.34	276	6	0.00	ng/ul#	1

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

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