

Data Path : Z:\HPCHEM1\BNA_M\Data\BM112816\
 Data File : BM008088.D
 Acq On : 28 Nov 2016 16:08
 Operator : UM/SJ
 Sample : H5695-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHST3

Quant Time: Nov 28 16:40:04 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 25 01:27:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	635	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.49	136	2173	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.35	164	1287	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	2424	0.40	ng/ul	-0.02
16) Chrysene-d12	21.29	240	2602	0.40	ng/ul	0.00
20) Perylene-d12	23.53	264	2646	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	420	0.14	ng/ul	-0.01
14) Fluoranthene-d10	19.13	212	2775	0.40	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	147	0.02	ng/ul#	11
5) 2-Methylnaphthalene	12.16	142	96	0.02	ng/ul	96
7) Acenaphthylene	14.06	152	151	0.03	ng/ul#	20
11) Pentachlorophenol	16.76	266	19	0.02	ng/ul#	70
12) Phenanthrene	17.12	178	1332	0.18	ng/ul#	87
13) Anthracene	17.23	178	379	0.05	ng/ul#	46
15) Fluoranthene	19.16	202	3772	0.39	ng/ul	90
17) Pyrene	19.52	202	3918	0.43	ng/ul#	90
18) Benzo(a)anthracene	21.30	228	10389	1.30	ng/ul#	67
19) Chrysene	21.30	228	10378	1.04	ng/ul#	69
21) Benzo(b)fluoranthene	22.86	252	7096	0.67	ng/ul#	24
22) Benzo(k)fluoranthene	22.86	252	7108	0.65	ng/ul#	20
23) Benzo(a)pyrene	23.43	252	2845	0.29	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.80	276	1702	0.14	ng/ul#	68
25) Dibenzo(a,h)anthracene	25.79	278	1519	0.16	ng/ul#	1
26) Benzo(g,h,i)perylene	26.51	276	9470	0.87	ng/ul#	64

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

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