

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032416\
 Data File : BM004766.D
 Acq On : 25 Mar 2016 01:02
 Operator : UM/SJ
 Sample : H1909-12
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4T46

Quant Time: Mar 25 03:06:55 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 25 03:18:48 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	74582	5.00	ng/ul	0.00
4) Naphthalene-d8	10.62	136	301093	5.00	ng/ul	0.00
8) Acenaphthene-d10	14.47	164	155655	5.00	ng/ul	0.00
12) Phenanthrene-d10	17.21	188	324348	5.00	ng/ul	0.00
18) Chrysene-d12	21.40	240	257648	5.00	ng/ul	-0.01
22) Perylene-d12	23.70	264	243038	5.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 1,4-Dioxane-d8	3.32	96	4037	0.11	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.22	152	1535	0.00	ng/ul	0.00
16) Fluoranthene-d10	19.24	212	3236	0.00	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
5) Naphthalene	10.67	128	2893	0.00	ng/ul#	91
9) Acenaphthylene	14.18	152	3357	0.01	ng/ul#	96
14) Phenanthrene	17.24	178	13786	0.02	ng/ul#	91
15) Anthracene	17.33	178	4342	0.01	ng/ul#	88
17) Fluoranthene	19.27	202	44981	0.05	ng/ul	95
19) Pyrene	19.63	202	39861	0.04	ng/ul	97
20) Benzo(a)anthracene	21.38	228	21047	0.03	ng/ul#	88
21) Chrysene	21.43	228	21353	0.03	ng/ul#	89
25) Benzo(a)pyrene	23.59	252	18654	0.03	ng/ul#	76
26) Indeno(1,2,3-cd)pyrene	26.02	276	11824	0.01	ng/ul#	89
28) Benzo(g,h,i)perylene	26.74	276	10697	0.01	ng/ul#	75

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

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