

Data Path : Z:\HPCHEM1\BNA M\DATA\BM032516\
 Data File : BM004795.D
 Acq On : 26 Mar 2016 02:22
 Operator : UM/SJ
 Sample : H1909-12
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4T46

Quant Time: Apr 05 19:44:04 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM032516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 05 18:29:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5	0.40	ng/ul	0.00
4) Naphthalene-d8	10.62	136	163	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.47	164	519	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.30	188	635742	0.40	ng/ul	0.09
18) Chrysene-d12	21.54	240	77	0.40	ng/ul	0.15
22) Perylene-d12	23.94	264	503	0.40	ng/ul	0.25
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.32	96	6998	3105.84	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.20	152	3911	17.21	ng/ul	0.00
16) Fluoranthene-d10	19.23	212	9517	0.01	ng/ul	-0.01
Target Compounds			Ovalue			
5) Naphthalene	10.67	128	7024	17.84	ng/ul	97
7) 2-Methylnaphthalene	12.28	142	1060	3.76	ng/ul	100
9) Acenaphthylene	14.18	152	9158	3.84	ng/ul#	94
10) Acenaphthene	14.52	153	796	0.47	ng/ul#	87
11) Fluorene	15.51	166	1962	0.97	ng/ul#	89
14) Phenanthrene	17.24	178	38235	0.02	ng/ul	97
19) Pyrene	19.63	202	114078	432.64	ng/ul	95
20) Benzo(a)anthracene	21.54	228	7140	30.17	ng/ul#	54
21) Chrysene	21.42	228	65429	286.15	ng/ul	95
23) Benzo(b)fluoranthene	23.16	252	14759	8.71	ng/ul	82
24) Benzo(k)fluoranthene	23.37	252	1130	0.72	ng/ul#	1
25) Benzo(a)pyrene	23.72	252	14575	9.41	ng/ul#	86
26) Indeno(1,2,3-cd)pyrene	26.27	276	1692	1.18	ng/ul#	33
27) Dibenzo(a,h)anthracene	26.27	278	6500	5.76	ng/ul#	83
28) Benzo(a,h,i)perylene	27.08	276	7111	5.99	ng/ul#	80

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

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