

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092216\
 Data File : BM007508.D
 Acq On : 22 Sep 2016 13:07
 Operator : UM/SJ
 Sample : H4830-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4VH7

Quant Time: Sep 22 16:27:59 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM092116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 22 01:26:49 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	1323	0.40	ng/ul	0.02
2) Naphthalene-d8	10.58	136	4877	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.42	164	3410	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.16	188	7788	0.40	ng/ul	0.00
16) Chrysene-d12	21.35	240	8482	0.40	ng/ul	0.00
20) Perylene-d12	23.62	264	7711	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.19	152	2137	0.38	ng/ul	0.02
14) Fluoranthene-d10	19.19	212	7666	0.42	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.63	128	11255	0.99	ng/ul	98
7) Acenaphthylene	14.14	152	11017	0.76	ng/ul	98
8) Acenaphthene	14.49	153	8640	0.77	ng/ul	97
9) Fluorene	15.48	166	12646	0.92	ng/ul	96
11) Pentachlorophenol	16.83	266	6136	2.89	ng/ul	98
12) Phenanthrene	17.21	178	17587	0.83	ng/ul	99
13) Anthracene	17.30	178	10172	0.53	ng/ul	99
17) Pyrene	19.58	202	48713	1.74	ng/ul	93
18) Benzo(a)anthracene	21.32	228	21374	0.85	ng/ul	94
19) Chrysene	21.38	228	31173	1.16	ng/ul	98
22) Benzo(k)fluoranthene	22.97	252	14322	0.55	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.90	276	20426	0.83	ng/ul	94
26) Benzo(g,h,i)perylene	26.60	276	23511	1.06	ng/ul	97

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

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