

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096187.D
 Acq On : 26 Apr 2018 19:56
 Operator : SJ/JU
 Sample : J2449-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Apr 27 05:40:18 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 27 04:45:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	722	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2957	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1717	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	3869m	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	5291m	0.40	ng/ul	0.00
20) Perylene-d12	24.36	264	3612m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	1571	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	4603m	0.26	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.23	128	530	0.064	ng/ul#	68
5) 2-Methylnaphthalene	12.81	142	333	0.060	ng/ul	96
8) Acenaphthene	15.01	153	222	0.035	ng/ul#	81
9) Fluorene	15.98	166	334	0.040	ng/ul#	81
12) Phenanthrene	17.70	178	6445m	0.566	ng/ul	
13) Anthracene	17.79	178	1208m	0.106	ng/ul	
15) Fluoranthene	19.67	202	14625m	0.840	ng/ul	
17) Pyrene	20.02	202	9328m	0.879	ng/ul	
18) Benzo(a)anthracene	21.73	228	4810m	0.300	ng/ul	
19) Chrysene	21.79	228	6317m	0.429	ng/ul	
21) Benzo(b)fluoranthene	23.55	252	8300m	0.607	ng/ul	
22) Benzo(k)fluoranthene	23.60	252	2619m	0.200	ng/ul	
23) Benzo(a)pyrene	24.24	252	5301m	0.446	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.14	276	3801m	0.237	ng/ul	
25) Dibenzo(a,h)anthracene	27.14	278	1135m	0.085	ng/ul	
26) Benzo(a,h,i)perylene	28.00	276	3197m	0.229	ng/ul	

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

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