

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095892.D
 Acq On : 29 Mar 2018 00:17
 Operator : SJ/JU
 Sample : J2060-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Mar 29 06:31:02 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 29 04:43:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.41	152	1722	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	4805	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	3304	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	7706m	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7707m	0.40	ng/ul	0.00
20) Perylene-d12	24.41	264	7704m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	2355	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	7762m	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.28	128	11184	0.834	ng/ul#	88
5) 2-Methylnaphthalene	12.87	142	4805	0.515	ng/ul	99
7) Acenaphthylene	14.72	152	554	0.033	ng/ul#	25
8) Acenaphthene	15.06	153	31382	2.498	ng/ul	96
9) Fluorene	16.02	166	39741	2.915	ng/ul#	90
12) Phenanthrene	17.75	178	332308m	14.399	ng/ul	
13) Anthracene	17.83	178	100300m	4.648	ng/ul	
15) Fluoranthene	19.71	202	523843m	18.739	ng/ul	
17) Pyrene	20.06	202	491802	16.805	ng/ul#	79
18) Benzo(a)anthracene	21.77	228	263300m	10.588	ng/ul	
19) Chrysene	21.83	228	243062	9.600	ng/ul	98
21) Benzo(b)fluoranthene	23.60	252	270642m	10.142	ng/ul	
22) Benzo(k)fluoranthene	23.65	252	92652m	3.652	ng/ul	
23) Benzo(a)pyrene	24.30	252	189143m	7.763	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.22	276	104003m	3.850	ng/ul	
25) Dibenzo(a,h)anthracene	27.23	278	33559m	1.554	ng/ul	
26) Benzo(a,h,i)perylene	28.09	276	91444m	3.940	ng/ul	

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

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