

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095893.D
 Acq On : 29 Mar 2018 00:53
 Operator : SJ/JU
 Sample : J2060-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Mar 29 06:36:10 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	1665	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	4763	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	3026	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	7338m	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7611m	0.40	ng/ul	0.00
20) Perylene-d12	24.41	264	8148	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	1597	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	4730m	0.20	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.28	128	739	0.056	ng/ul#	64
5) 2-Methylnaphthalene	12.87	142	402	0.043	ng/ul	98
8) Acenaphthene	15.06	153	274	0.024	ng/ul#	69
9) Fluorene	16.02	166	328	0.026	ng/ul#	76
12) Phenanthrene	17.74	178	2111m	0.096	ng/ul	
13) Anthracene	17.83	178	791m	0.038	ng/ul	
15) Fluoranthene	19.71	202	4002m	0.150	ng/ul	
17) Pyrene	20.06	202	5612	0.194	ng/ul#	84
18) Benzo(a)anthracene	21.77	228	2553m	0.104	ng/ul	
19) Chrysene	21.82	228	2911m	0.116	ng/ul	
21) Benzo(b)fluoranthene	23.60	252	3658m	0.130	ng/ul	
22) Benzo(k)fluoranthene	23.64	252	1409m	0.053	ng/ul	
23) Benzo(a)pyrene	24.30	252	2748	0.107	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	27.21	276	1941	0.068	ng/ul#	76
26) Benzo(g,h,i)perylene	28.08	276	2086	0.085	ng/ul#	91

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

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