

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE090718\
 Data File : BE097420.D
 Acq On : 7 Sep 2018 12:05
 Operator : SJ/JU
 Sample : J4688-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z17

Manual Integrations
 APPROVED

Sohil
 9/11/2018 2:42:26 PM

Quant Time: Sep 08 00:10:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 07 23:49:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	972	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	5009	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3098	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	8323m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9850m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	8815m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1008	0.13	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	3660m	0.13	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.17	128	338	0.028	ng/ul#	67
5) 2-Methylnaphthalene	11.80	142	272	0.035	ng/ul	96
12) Phenanthrene	16.80	178	1914m	0.089	ng/ul	
15) Fluoranthene	18.83	202	3211m	0.117	ng/ul	
17) Pyrene	19.19	202	3057m	0.115	ng/ul	
18) Benzo(a)anthracene	20.95	228	1572	0.062	ng/ul#	88
19) Chrysene	21.00	228	3036	0.104	ng/ul#	91
21) Benzo(b)fluoranthene	22.53	252	2854m	0.121	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	1098m	0.042	ng/ul	
23) Benzo(a)pyrene	23.10	252	1496	0.069	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.49	276	1267	0.045	ng/ul#	84
26) Benzo(g,h,i)perylene	26.19	276	1264	0.052	ng/ul#	92

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

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