

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102218\
 Data File : BE098007.D
 Acq On : 22 Oct 2018 13:11
 Operator : SJ/JU
 Sample : J5567-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AH4

Manual Integrations
 APPROVED

Sohil
 10/23/2018 5:34:09 PM

Quant Time: Oct 23 02:04:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 01:24:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	2778	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	13557	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	9705	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	22309m	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	23694m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	21809m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	7861	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	22185m	0.31	ng/ul	0.00
Target Compounds				Qvalue		
15) Fluoranthene	19.35	202	2228m	0.032	ng/ul	
17) Pyrene	19.71	202	1593	0.027	ng/ul#	42
18) Benzo(a)anthracene	21.45	228	2942m	0.045	ng/ul	
19) Chrysene	21.51	228	2125m	0.033	ng/ul	
21) Benzo(b)fluoranthene	23.25	252	7521m	0.133	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	1720m	0.027	ng/ul	
23) Benzo(a)pyrene	23.91	252	1866m	0.036	ng/ul	

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

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