

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102218\
 Data File : BE098018.D
 Acq On : 22 Oct 2018 20:34
 Operator : SJ/JU
 Sample : J5567-12
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AJ0

Manual Integrations
 APPROVED

Sohil
 10/23/2018 5:34:38 PM

Quant Time: Oct 23 03:05:25 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 02:34:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.88	152	1625	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	8070	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	5865	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	13335m	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	14435m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	13974m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	3917	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	11173m	0.26	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.34	178	1400m	0.044	ng/ul	
15) Fluoranthene	19.36	202	2915m	0.069	ng/ul	
17) Pyrene	19.72	202	1787m	0.050	ng/ul	
18) Benzo(a)anthracene	21.46	228	2350m	0.059	ng/ul	
19) Chrysene	21.51	228	1704m	0.044	ng/ul	
21) Benzo(b)fluoranthene	23.24	252	4642m	0.129	ng/ul	
22) Benzo(k)fluoranthene	23.28	252	1111m	0.027	ng/ul	
23) Benzo(a)pyrene	23.92	252	1327m	0.040	ng/ul	

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

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