

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102218\
 Data File : BE098012.D
 Acq On : 22 Oct 2018 16:16
 Operator : SJ/JU
 Sample : J5567-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AH8

Manual Integrations
 APPROVED

Sohil
 10/23/2018 5:34:26 PM

Quant Time: Oct 23 02:28:48 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	1855	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	9011	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	6549	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	14845m	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	15833m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	14122m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6016	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	16647m	0.34	ng/ul	0.00
Target Compounds					Qvalue	
12) Phenanthrene	17.34	178	764m	0.022	ng/ul	
15) Fluoranthene	19.36	202	4500m	0.096	ng/ul	
17) Pyrene	19.72	202	3040	0.078	ng/ul#	62
18) Benzo(a)anthracene	21.46	228	4519	0.103	ng/ul#	85
19) Chrysene	21.51	228	3897	0.092	ng/ul#	80
21) Benzo(b)fluoranthene	23.25	252	12041m	0.330	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	2427m	0.059	ng/ul	
23) Benzo(a)pyrene	23.92	252	3559	0.107	ng/ul#	48
24) Indeno(1,2,3-cd)pyrene	26.74	276	1335	0.035	ng/ul#	100
26) Benzo(g,h,i)perylene	27.56	276	1485	0.041	ng/ul#	36

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

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