

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE102218\
 Data File : BE098009.D
 Acq On : 22 Oct 2018 14:26
 Operator : SJ/JU
 Sample : J5567-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AH2

Manual Integrations
 APPROVED

Sohil
 10/23/2018 5:34:13 PM

Quant Time: Oct 23 02:10:44 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	2865	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	13693	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	9149	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	21113	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	26349	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	25312	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	4078	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	10978	0.16	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.34	178	1238	0.025	ng/ul#	78
13) Anthracene	17.43	178	1067	0.025	ng/ul#	54
15) Fluoranthene	19.35	202	8089	0.122	ng/ul#	59
17) Pyrene	19.72	202	5903m	0.091	ng/ul	
18) Benzo(a)anthracene	21.46	228	8896	0.122	ng/ul#	82
19) Chrysene	21.51	228	7563	0.107	ng/ul#	82
21) Benzo(b)fluoranthene	23.25	252	20927m	0.320	ng/ul	
22) Benzo(k)fluoranthene	23.30	252	4234m	0.057	ng/ul	
23) Benzo(a)pyrene	23.91	252	5503	0.092	ng/ul#	55
24) Indeno(1,2,3-cd)pyrene	26.73	276	2095	0.031	ng/ul#	100
26) Benzo(g,h,i)perylene	27.56	276	2528	0.039	ng/ul#	55

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

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