

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080818\
 Data File : BE097233.D
 Acq On : 8 Aug 2018 13:47
 Operator : SJ/JU
 Sample : J4268-13
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AC3

Manual Integrations
 APPROVED

Sohil
 8/9/2018 4:43:02 PM

Quant Time: Aug 09 00:30:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 08 23:25:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1255	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6640	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4027	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	9399m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	9567m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	9576m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1999	0.19	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	5835m	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	3386	0.211	ng/ul#	95
5) 2-Methylnaphthalene	11.80	142	1811	0.170	ng/ul	98
7) Acenaphthylene	13.74	152	7220	0.415	ng/ul	97
8) Acenaphthene	14.08	153	3143	0.250	ng/ul#	93
9) Fluorene	15.08	166	4708	0.325	ng/ul	93
12) Phenanthrene	16.81	178	144987m	6.051	ng/ul	
13) Anthracene	16.90	178	21875m	0.979	ng/ul	
15) Fluoranthene	18.84	202	487348m	15.388	ng/ul	
17) Pyrene	19.21	202	436060	17.010	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	233018m	8.524	ng/ul	
19) Chrysene	21.01	228	279425	9.919	ng/ul	96
21) Benzo(b)fluoranthene	22.54	252	396895m	15.207	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	130211m	4.627	ng/ul	
23) Benzo(a)pyrene	23.12	252	254921m	9.761	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.51	276	205262	6.280	ng/ul#	70
25) Dibenzo(a,h)anthracene	25.52	278	46555m	1.741	ng/ul	
26) Benzo(a,h,i)perylene	26.21	276	200753	7.107	ng/ul#	90

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

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