

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080818\
 Data File : BE097240.D
 Acq On : 8 Aug 2018 18:13
 Operator : SJ/JU
 Sample : J4268-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA6

Manual Integrations
 APPROVED

Sohil
 8/9/2018 4:43:21 PM

Quant Time: Aug 09 01:34:30 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	1147	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6064	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3674	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	8579m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	8752m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	7996m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1878	0.19	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	5430m	0.18	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	1120	0.076	ng/ul#	89
5) 2-Methylnaphthalene	11.80	142	256	0.026	ng/ul	97
7) Acenaphthylene	13.73	152	602	0.038	ng/ul#	77
8) Acenaphthene	14.08	153	601	0.052	ng/ul#	92
9) Fluorene	15.08	166	793	0.060	ng/ul#	93
12) Phenanthrene	16.81	178	12112m	0.554	ng/ul	
13) Anthracene	16.90	178	2239m	0.110	ng/ul	
15) Fluoranthene	18.84	202	31772m	1.099	ng/ul	
17) Pyrene	19.20	202	31446	1.341	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	17002m	0.680	ng/ul	
19) Chrysene	21.01	228	17418	0.676	ng/ul	98
21) Benzo(b)fluoranthene	22.53	252	23462m	1.077	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	7005m	0.298	ng/ul	
23) Benzo(a)pyrene	23.11	252	16265m	0.746	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	10792m	0.395	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	2595m	0.116	ng/ul	
26) Benzo(a,h,i)perylene	26.21	276	10838m	0.460	ng/ul	

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

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