

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080818\
 Data File : BE097239.D
 Acq On : 8 Aug 2018 17:34
 Operator : SJ/JU
 Sample : J4268-12
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AA2

Manual Integrations
 APPROVED

Sohil
 8/9/2018 4:43:19 PM

Quant Time: Aug 09 01:19:15 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	1154	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6170	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3704	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	8570m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	8715m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8229m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1577	0.16	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	4397m	0.15	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.08	153	662	0.057	ng/ul#	92
9) Fluorene	15.08	166	1183	0.089	ng/ul	88
12) Phenanthrene	16.81	178	21146m	0.968	ng/ul	
13) Anthracene	16.90	178	4081m	0.200	ng/ul	
15) Fluoranthene	18.84	202	44893m	1.555	ng/ul	
17) Pyrene	19.20	202	37906	1.623	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	19360m	0.777	ng/ul	
19) Chrysene	21.01	228	19807m	0.772	ng/ul	
21) Benzo(b)fluoranthene	22.54	252	26229m	1.169	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	7890m	0.326	ng/ul	
23) Benzo(a)pyrene	23.11	252	17302m	0.771	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	11752m	0.418	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	2817m	0.123	ng/ul	
26) Benzo(g,h,i)perylene	26.21	276	11165m	0.460	ng/ul	

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

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