

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080318\
 Data File : BE097174.D
 Acq On : 3 Aug 2018 10:32
 Operator : SJ/JU
 Sample : J4172-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0CZ8

Manual Integrations
 APPROVED

Sohil
 8/6/2018 4:32:18 PM

Quant Time: Aug 03 12:01:06 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 : Fri Aug 03 11:26:05 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1490	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	7816	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	5006	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11337m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	12310m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	11422m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	3698	0.29	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	10627m	0.27	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	1929	0.102	ng/ul#	94
5) 2-Methylnaphthalene	11.80	142	1392	0.111	ng/ul	98
7) Acenaphthylene	13.74	152	978	0.045	ng/ul#	73
8) Acenaphthene	14.08	153	19204	1.227	ng/ul	96
9) Fluorene	15.08	166	12356	0.687	ng/ul	92
12) Phenanthrene	16.81	178	235607m	8.152	ng/ul	
13) Anthracene	16.90	178	67232m	2.494	ng/ul	
15) Fluoranthene	18.84	202	775921m	20.311	ng/ul	
17) Pyrene	19.21	202	736006	22.313	ng/ul#	78
18) Benzo(a)anthracene	20.96	228	450432m	12.805	ng/ul	
19) Chrysene	21.01	228	423147	11.674	ng/ul	98
21) Benzo(b)fluoranthene	22.54	252	636828m	20.456	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	219350m	6.535	ng/ul	
23) Benzo(a)pyrene	23.12	252	472794m	15.177	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.50	276	318714m	8.175	ng/ul	
25) Dibenzo(a,h)anthracene	25.52	278	75878m	2.378	ng/ul	
26) Benzo(a,h,i)perylene	26.21	276	281103m	8.343	ng/ul	

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

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