

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101218\
 Data File : BE097889.D
 Acq On : 12 Oct 2018 21:16
 Operator : SJ/JU
 Sample : J5183-08DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC47DL

Manual Integrations
 APPROVED

Sohil
 10/15/2018 3:55:49 PM

Quant Time: Oct 13 03:08:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 13 02:57:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1903	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	9698	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.56	164	6303	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.32	188	15299m	0.40	ng/ul	0.00
16) Chrysene-d12	21.50	240	15857m	0.40	ng/ul	0.00
20) Perylene-d12	24.07	264	15795m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1065	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.35	212	3391m	0.07	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	723	0.031	ng/ul#	70
5) 2-Methylnaphthalene	12.37	142	552	0.032	ng/ul	90
8) Acenaphthene	14.62	153	1282	0.069	ng/ul	96
9) Fluorene	15.61	166	1537	0.064	ng/ul#	96
12) Phenanthrene	17.35	178	29120	0.804	ng/ul	98
13) Anthracene	17.44	178	5539m	0.165	ng/ul	
15) Fluoranthene	19.37	202	47516m	1.022	ng/ul	
17) Pyrene	19.74	202	59724	1.264	ng/ul	99
18) Benzo(a)anthracene	21.48	228	24514m	0.511	ng/ul	
19) Chrysene	21.54	228	56851m	1.320	ng/ul	
21) Benzo(b)fluoranthene	23.29	252	59718m	1.415	ng/ul	
22) Benzo(k)fluoranthene	23.33	252	18316m	0.405	ng/ul	
23) Benzo(a)pyrene	23.95	252	26660m	0.643	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.78	276	31012m	0.659	ng/ul	
25) Dibenzo(a,h)anthracene	26.80	278	7705m	0.198	ng/ul	
26) Benzo(g,h,i)perylene	27.62	276	34289m	0.836	ng/ul	

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

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