

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102617\
 Data File : BE094588.D
 Acq On : 27 Oct 2017 10:19
 Operator : SJ/JU
 Sample : I5786-06 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-8(5-6.25)_101017

Manual Integrations
 APPROVED

Sohil
 10/27/2017 5:36:37 PM

Quant Time: Oct 27 13:27:25 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 27 04:05:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1836	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10306	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	5892	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	12750	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	12144	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	12677	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	944	0.06	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	2275	0.06	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	2102	0.083	ng/ul#	90
5) 2-Methylnaphthalene	12.35	142	1547	0.088	ng/ul	100
7) Acenaphthylene	14.24	152	3748	0.141	ng/ul	95
8) Acenaphthene	14.58	153	1015	0.053	ng/ul#	91
9) Fluorene	15.57	166	2585	0.120	ng/ul	90
12) Phenanthrene	17.29	178	33229m	0.987	ng/ul	
13) Anthracene	17.40	178	7201	0.217	ng/ul#	93
15) Fluoranthene	19.32	202	47558	1.221	ng/ul	84
17) Pyrene	19.68	202	44625	1.189	ng/ul#	81
18) Benzo(a)anthracene	21.41	228	21558	0.621	ng/ul#	87
19) Chrysene	21.47	228	24134	0.673	ng/ul#	85
21) Benzo(b)fluoranthene	23.19	252	29758m	0.834	ng/ul	
22) Benzo(k)fluoranthene	23.23	252	9551m	0.250	ng/ul	
23) Benzo(a)pyrene	23.85	252	21293	0.597	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	26.67	276	13386	0.363	ng/ul#	75
25) Dibenzo(a,h)anthracene	26.66	278	3408	0.110	ng/ul#	1
26) Benzo(g,h,i)perylene	27.52	276	11422	0.386	ng/ul#	63

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

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