

Data Path : Z:\HPCHEM1\BNA E\DATA\BE112217\
 Data File : BE094842.D
 Acq On : 22 Nov 2017 21:17
 Operator : SJ/JU
 Sample : I6496-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
ClientSampleId :
 C0AA1

Manual Integrations
APPROVED
 Sohil
 11/27/2017 3:25:41 PM

Quant Time: Nov 23 03:05:02 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE111417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 23 02:50:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	
1) 1,4-Dichlorobenzene-d4	7.89	152	691	0.40	ng/ul	-0.04	
2) Naphthalene-d8	10.67	136	4662	0.40	ng/ul	-0.04	
6) Acenaphthene-d10	14.50	164	3365	0.40	ng/ul	-0.01	
10) Phenanthrene-d10	17.24	188	8284m	0.40	ng/ul	-0.02	
16) Chrysene-d12	21.42	240	8449	0.40	ng/ul	0.00	
20) Perylene-d12	23.96	264	9439	0.40	ng/ul	-0.01	
System Monitoring Compounds							
4) 2-Methylnaphthalene-d10	12.29	152	2197	0.29	ng/ul	0.00	
14) Fluoranthene-d10	19.27	212	7478	0.26	ng/ul	-0.01	
Target Compounds							
							Ovalue
3) Naphthalene	10.72	128	449	0.040	ng/ul#		64
5) 2-Methylnaphthalene	12.37	142	587	0.076	ng/ul		97
12) Phenanthrene	17.27	178	2946m	0.132	ng/ul		
13) Anthracene	17.36	178	585m	0.026	ng/ul		
15) Fluoranthene	19.30	202	5035	0.172	ng/ul		84
17) Pyrene	19.66	202	5495	0.208	ng/ul#		85
18) Benzo(a)anthracene	21.41	228	2431	0.095	ng/ul#		90
19) Chrysene	21.46	228	3411	0.133	ng/ul#		91
21) Benzo(b)fluoranthene	23.17	252	3948m	0.140	ng/ul		
22) Benzo(k)fluoranthene	23.21	252	1247m	0.042	ng/ul		
23) Benzo(a)pyrene	23.83	252	2259	0.081	ng/ul#		92
24) Indeno(1,2,3-cd)pyrene	26.68	276	1663	0.055	ng/ul#		77
26) Benzo(g,h,i)perylene	27.53	276	1420	0.056	ng/ul#		71

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

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