

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102617\
 Data File : BE094569.D
 Acq On : 26 Oct 2017 19:15
 Operator : SJ/JU
 Sample : I5719-16 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-26(0.5-1.5)-100517

Manual Integrations
 APPROVED

Sohil
 10/27/2017 5:35:34 PM

Quant Time: Oct 27 08:45:33 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	7.90	152	1014	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	5672	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.51	164	3436	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	7876	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	8203	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	8172	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	433	0.05	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	1229	0.05	ng/ul	0.00
Target Compounds						
					Qvalue	
3) Naphthalene	10.74	128	2699	0.195	ng/ul#	94
5) 2-Methylnaphthalene	12.35	142	1712	0.176	ng/ul	97
9) Fluorene	15.58	166	376	0.030	ng/ul#	1
12) Phenanthrene	17.29	178	14078	0.677	ng/ul#	91
13) Anthracene	17.38	178	796	0.039	ng/ul#	71
15) Fluoranthene	19.31	202	8782	0.365	ng/ul	83
17) Pyrene	19.67	202	9774	0.385	ng/ul#	85
18) Benzo(a)anthracene	21.40	228	3655	0.156	ng/ul#	62
19) Chrysene	21.46	228	11696	0.483	ng/ul#	87
21) Benzo(b)fluoranthene	23.18	252	6823m	0.297	ng/ul	
22) Benzo(k)fluoranthene	23.21	252	1834m	0.074	ng/ul	
23) Benzo(a)pyrene	23.85	252	3130	0.136	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.65	276	2573	0.108	ng/ul#	85
25) Dibenzo(a,h)anthracene	26.64	278	1221	0.061	ng/ul#	1
26) Benzo(g,h,i)perylene	27.50	276	3172	0.166	ng/ul#	1

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

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