

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110617\
 Data File : BE094762.D
 Acq On : 6 Nov 2017 20:26
 Operator : SJ/JU
 Sample : I6096-03 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B-52(0-1)-102317

Manual Integrations
 APPROVED

Sohil
 11/7/2017 5:41:53 PM

Quant Time: Nov 07 01:38:32 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 07 01:37:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1459	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.69	136	8275	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.51	164	4571	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10416	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	10015	0.40	ng/ul	0.00
20) Perylene-d12	23.97	264	10247	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	453	0.03	ng/ul	-0.02
14) Fluoranthene-d10	19.28	212	826	0.03	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.73	128	44756	2.212	ng/ul#	93
5) 2-Methylnaphthalene	12.34	142	7200	0.509	ng/ul	97
7) Acenaphthylene	14.24	152	1732	0.084	ng/ul#	81
8) Acenaphthene	14.57	153	2704	0.182	ng/ul	96
9) Fluorene	15.56	166	4531	0.272	ng/ul	92
12) Phenanthrene	17.29	178	84264	3.063	ng/ul#	90
13) Anthracene	17.38	178	11290	0.416	ng/ul#	80
15) Fluoranthene	19.31	202	122068	3.838	ng/ul	81
17) Pyrene	19.67	202	119046	3.845	ng/ul#	83
18) Benzo(a)anthracene	21.41	228	76941	2.687	ng/ul#	89
19) Chrysene	21.47	228	88079	2.976	ng/ul	96
21) Benzo(b)fluoranthene	23.19	252	102167m	3.544	ng/ul	
22) Benzo(k)fluoranthene	23.24	252	29751m	0.963	ng/ul	
23) Benzo(a)pyrene	23.86	252	64521	2.239	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.69	276	38511	1.292	ng/ul#	73
25) Dibenzo(a,h)anthracene	26.68	278	9421	0.375	ng/ul#	58
26) Benzo(g,h,i)perylene	27.54	276	32364	1.352	ng/ul#	92

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

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