

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE110617\
 Data File : BE094766.D
 Acq On : 6 Nov 2017 22:49
 Operator : SJ/JU
 Sample : I6096-07 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 DUP-7

Manual Integrations
 APPROVED

Sohil
 11/7/2017 5:42:33 PM

Quant Time: Nov 07 01:52:26 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 01:23:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1440	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.69	136	8486	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.52	164	4767	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	10603	0.40	ng/ul	-0.02
16) Chrysene-d12	21.43	240	10729	0.40	ng/ul	0.00
20) Perylene-d12	23.98	264	10893	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	484	0.04	ng/ul	-0.01
14) Fluoranthene-d10	19.29	212	1113	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	5556	0.268	ng/ul#	94
5) 2-Methylnaphthalene	12.36	142	1171	0.081	ng/ul	97
7) Acenaphthylene	14.24	152	17209	0.798	ng/ul	96
8) Acenaphthene	14.58	153	1149	0.074	ng/ul#	92
9) Fluorene	15.57	166	1708	0.098	ng/ul	92
12) Phenanthrene	17.29	178	44587m	1.592	ng/ul	
13) Anthracene	17.39	178	12778	0.462	ng/ul	97
15) Fluoranthene	19.31	202	134808	4.164	ng/ul	84
17) Pyrene	19.68	202	151524	4.568	ng/ul#	78
18) Benzo(a)anthracene	21.41	228	91466	2.982	ng/ul#	89
19) Chrysene	21.47	228	114859	3.623	ng/ul	95
21) Benzo(b)fluoranthene	23.19	252	236015m	7.702	ng/ul	
22) Benzo(k)fluoranthene	23.22	252	79986m	2.436	ng/ul	
23) Benzo(a)pyrene	23.86	252	161792	5.282	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	26.69	276	100804	3.181	ng/ul#	74
25) Dibenzo(a,h)anthracene	26.68	278	24135	0.904	ng/ul#	90
26) Benzo(g,h,i)perylene	27.53	276	80733	3.172	ng/ul	95

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

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