

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102717\
 Data File : BE094613.D
 Acq On : 28 Oct 2017 1:35
 Operator : SJ/JU
 Sample : I5842-16
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AB4

Manual Integrations
 APPROVED

Sohil
 10/30/2017 6:37:17 PM

Quant Time: Oct 30 02:38:01 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 30 02:20:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1094	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	6477	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.52	164	3980	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	8636m	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	7983	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	7904m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4186	0.40	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	10569	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	820	0.052	ng/ul#	81
5) 2-Methylnaphthalene	12.35	142	663	0.060	ng/ul	99
8) Acenaphthene	14.58	153	656	0.051	ng/ul	96
9) Fluorene	15.57	166	599	0.041	ng/ul#	94
12) Phenanthrene	17.29	178	18192	0.798	ng/ul#	89
13) Anthracene	17.38	178	2907	0.129	ng/ul#	94
15) Fluoranthene	19.31	202	54141	2.053	ng/ul	84
17) Pyrene	19.67	202	48393	1.961	ng/ul#	80
18) Benzo(a)anthracene	21.41	228	20743	0.909	ng/ul#	88
19) Chrysene	21.46	228	27915	1.183	ng/ul	98
21) Benzo(b)fluoranthene	23.18	252	37590	1.691	ng/ul	88
22) Benzo(k)fluoranthene	23.23	252	15612m	0.655	ng/ul	
23) Benzo(a)pyrene	23.85	252	22745m	1.023	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.67	276	15476m	0.673	ng/ul	
25) Dibenzo(a,h)anthracene	26.66	278	3264m	0.168	ng/ul	
26) Benzo(a,h,i)perylene	27.51	276	11334m	0.614	ng/ul	

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

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