

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100917\
 Data File : BE094245.D
 Acq On : 10 Oct 2017 6:05
 Operator : SJ/JU
 Sample : I5533-07
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 DUP-1

Manual Integrations
 APPROVED

Sohil
 10/11/2017 9:24:22 AM

Quant Time: Oct 10 12:23:01 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 09 04:05:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	2346	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	11916	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	7160	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.26	188	14817	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	13653	0.40	ng/ul	0.00
20) Perylene-d12	23.93	264	13389m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	4314	0.23	ng/ul	-0.01
14) Fluoranthene-d10	19.28	212	9856	0.20	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	26639	0.913	ng/ul#	91
5) 2-Methylnaphthalene	12.35	142	8277	0.437	ng/ul	98
7) Acenaphthylene	14.24	152	22417	0.803	ng/ul	97
8) Acenaphthene	14.57	153	37357	1.626	ng/ul	96
9) Fluorene	15.56	166	39873	1.482	ng/ul	90
12) Phenanthrene	17.29	178	873416	22.557	ng/ul#	89
13) Anthracene	17.38	178	220972	5.868	ng/ul#	89
15) Fluoranthene	19.31	202	1536532	30.201	ng/ul	84
17) Pyrene	19.67	202	1442462	37.972	ng/ul#	86
18) Benzo(a)anthracene	21.40	228	758875	20.952	ng/ul#	86
19) Chrysene	21.45	228	678967	16.233	ng/ul	99
21) Benzo(b)fluoranthene	23.16	252	810255m	22.277	ng/ul	
22) Benzo(k)fluoranthene	23.20	252	270517m	6.435	ng/ul	
23) Benzo(a)pyrene	23.82	252	604156m	15.985	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.59	276	374841m	9.695	ng/ul	
25) Dibenzo(a,h)anthracene	26.59	278	103530m	3.193	ng/ul	
26) Benzo(a,h,i)perylene	27.42	276	308326m	9.246	ng/ul	

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

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