

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094273.D
 Acq On : 11 Oct 2017 4:42
 Operator : SJ/JU
 Sample : I5522-02
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AD3

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:20:33 PM

Quant Time: Oct 11 08:33:48 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 08:24:24 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	2183	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	10897	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	6386	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.24	188	12734m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.41	240	12887	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	14508m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.27	152	4068	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	9562	0.23	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	21662	0.812	ng/ul#	91
5) 2-Methylnaphthalene	12.35	142	7062	0.408	ng/ul	98
7) Acenaphthylene	14.24	152	5586	0.224	ng/ul	98
8) Acenaphthene	14.57	153	28509	1.391	ng/ul	97
9) Fluorene	15.56	166	35369	1.474	ng/ul#	93
12) Phenanthrene	17.29	178	640298	19.242	ng/ul#	93
13) Anthracene	17.38	178	146919	4.540	ng/ul#	91
15) Fluoranthene	19.30	202	1336654	30.570	ng/ul	85
17) Pyrene	19.66	202	1113467	31.054	ng/ul#	84
18) Benzo(a)anthracene	21.39	228	595830	17.428	ng/ul#	86
19) Chrysene	21.45	228	641834	16.257	ng/ul	99
21) Benzo(b)fluoranthene	23.15	252	716234	18.173	ng/ul	85
22) Benzo(k)fluoranthene	23.19	252	256021m	5.620	ng/ul	
23) Benzo(a)pyrene	23.81	252	511244m	12.484	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.58	276	340747m	8.134	ng/ul	
25) Dibenzo(a,h)anthracene	26.58	278	90941m	2.588	ng/ul	
26) Benzo(a,h,i)perylene	27.41	276	276377m	7.648	ng/ul	

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

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