

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093296.D
 Acq On : 21 Jun 2017 00:05
 Operator : SJ/JU
 Sample : I3627-16
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A89

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:19:05 PM

Quant Time: Jun 21 04:04:39 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 21 03:23:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	2103	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	10959	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	7580	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	17005	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	16403	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	13799m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	4717	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	12890	0.28	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.56	128	27033	0.994	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	16513	0.888	ng/ul	98
7) Acenaphthylene	14.09	152	5028	0.155	ng/ul#	94
8) Acenaphthene	14.42	153	45819	1.746	ng/ul	97
9) Fluorene	15.42	166	57416	1.866	ng/ul	94
12) Phenanthrene	17.14	178	521671	11.149	ng/ul#	87
13) Anthracene	17.22	178	51479	1.264	ng/ul#	88
15) Fluoranthene	19.15	202	305455	5.324	ng/ul	84
17) Pyrene	19.52	202	186273	3.851	ng/ul#	84
18) Benzo(a)anthracene	21.24	228	48019	1.097	ng/ul#	85
19) Chrysene	21.29	228	41637	0.917	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	25054	0.598	ng/ul	88
22) Benzo(k)fluoranthene	22.97	252	9318m	0.210	ng/ul	
23) Benzo(a)pyrene	23.56	252	12815m	0.324	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.18	276	5461m	0.119	ng/ul	
25) Dibenzo(a,h)anthracene	26.19	278	1818m	0.047	ng/ul	
26) Benzo(g,h,i)perylene	26.95	276	4286	0.109	ng/ul#	90

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

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