

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093288.D
 Acq On : 20 Jun 2017 19:10
 Operator : SJ/JU
 Sample : I3627-19
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A92

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:18:34 PM

Quant Time: Jun 21 03:18:34 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 01:49:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1701	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8733	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5593	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13418m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13073	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11370m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	3915	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	10655	0.29	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	372499	17.191	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	29490	1.990	ng/ul	98
7) Acenaphthylene	14.09	152	2505	0.104	ng/ul#	93
8) Acenaphthene	14.42	153	15976	0.825	ng/ul	98
9) Fluorene	15.42	166	28625	1.261	ng/ul#	94
12) Phenanthrene	17.14	178	121509	3.291	ng/ul#	88
13) Anthracene	17.22	178	10818	0.337	ng/ul#	92
15) Fluoranthene	19.15	202	22650	0.500	ng/ul	84
17) Pyrene	19.51	202	18296	0.475	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	3361	0.096	ng/ul#	86
19) Chrysene	21.30	228	3544	0.098	ng/ul	97
21) Benzo(b)fluoranthene	22.93	252	2505	0.073	ng/ul	87
22) Benzo(k)fluoranthene	22.98	252	1878m	0.051	ng/ul	
23) Benzo(a)pyrene	23.55	252	1878m	0.058	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.19	276	1835	0.049	ng/ul#	90
25) Dibenzo(a,h)anthracene	26.20	278	1386	0.044	ng/ul#	61
26) Benzo(g,h,i)perylene	26.95	276	1614	0.050	ng/ul#	77

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

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