

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093301.D
 Acq On : 21 Jun 2017 3:10
 Operator : SJ/JU
 Sample : I3599-05DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A81DL

Manual Integrations
 APPROVED

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

Quant Time: Jun 21 04:33:37 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	7.73	152	1994	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	10975	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	6674	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	15973	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	16413	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	14865	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	582	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	1485	0.03	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	964	0.035	ng/ul#	86
7) Acenaphthylene	14.08	152	5113	0.179	ng/ul	98
12) Phenanthrene	17.14	178	5066	0.115	ng/ul#	91
13) Anthracene	17.22	178	7006	0.183	ng/ul#	90
15) Fluoranthene	19.15	202	32053	0.595	ng/ul	83
17) Pyrene	19.51	202	38754	0.801	ng/ul#	85
18) Benzo(a)anthracene	21.24	228	23829	0.544	ng/ul#	85
19) Chrysene	21.29	228	27559	0.607	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	51732m	1.147	ng/ul	
22) Benzo(k)fluoranthene	22.97	252	16459m	0.344	ng/ul	
23) Benzo(a)pyrene	23.55	252	23707	0.556	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.17	276	16983	0.344	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	4331	0.104	ng/ul#	87
26) Benzo(g,h,i)perylene	26.95	276	13059	0.308	ng/ul	96

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

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