

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093286.D
 Acq On : 20 Jun 2017 17:55
 Operator : SJ/JU
 Sample : I3627-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A45

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:18:28 PM

Quant Time: Jun 21 03:13:34 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	1691	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	8827	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	5450	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	13027	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	12854	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	10029m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	2589	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	7215	0.20	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	472	0.022	ng/ul#	76
12) Phenanthrene	17.13	178	10293	0.287	ng/ul#	89
13) Anthracene	17.22	178	2178	0.070	ng/ul#	93
15) Fluoranthene	19.15	202	18339	0.417	ng/ul	83
17) Pyrene	19.51	202	16534	0.436	ng/ul#	88
18) Benzo(a)anthracene	21.24	228	3693	0.108	ng/ul#	85
19) Chrysene	21.30	228	3998	0.112	ng/ul	96
21) Benzo(b)fluoranthene	22.93	252	3768	0.124	ng/ul	85
22) Benzo(k)fluoranthene	22.98	252	1218m	0.038	ng/ul	
23) Benzo(a)pyrene	23.55	252	2192m	0.076	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.18	276	1375m	0.041	ng/ul	
26) Benzo(g,h,i)perylene	26.95	276	1061	0.037	ng/ul#	65

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

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