

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062017\
 Data File : BE093303.D
 Acq On : 21 Jun 2017 4:25
 Operator : SJ/JU
 Sample : I3599-06DL 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F5A87DL

Manual Integrations
 APPROVED

mohammad
 6/21/2017 12:19:27 PM

Quant Time: Jun 21 05:09:32 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	2023	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	10658	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	6299	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	15649	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	16586	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	14450	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	743	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	2087	0.05	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	24018	0.908	ng/ul#	91
5) 2-Methylnaphthalene	12.18	142	4945	0.273	ng/ul	99
7) Acenaphthylene	14.09	152	35718	1.321	ng/ul	96
8) Acenaphthene	14.42	153	7159	0.328	ng/ul	96
9) Fluorene	15.42	166	7317	0.286	ng/ul	96
12) Phenanthrene	17.14	178	77954	1.810	ng/ul#	90
13) Anthracene	17.22	178	590647	15.755	ng/ul#	88
15) Fluoranthene	19.16	202	1419499	26.883	ng/ul	83
17) Pyrene	19.52	202	1446524	29.573	ng/ul#	86
18) Benzo(a)anthracene	21.24	228	748096	16.909	ng/ul#	84
19) Chrysene	21.30	228	828744	18.058	ng/ul	98
21) Benzo(b)fluoranthene	22.94	252	952127m	21.712	ng/ul	
22) Benzo(k)fluoranthene	22.98	252	301880m	6.482	ng/ul	
23) Benzo(a)pyrene	23.56	252	436919	10.537	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.19	276	241304	5.026	ng/ul#	87
25) Dibenzo(a,h)anthracene	26.19	278	73082m	1.808	ng/ul	
26) Benzo(g,h,i)perylene	26.96	276	135491	3.291	ng/ul	96

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

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