

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081617\
 Data File : BE093839.D
 Acq On : 16 Aug 2017 14:08
 Operator : SJ/JU
 Sample : I4672-05
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AH1

Manual Integrations
 APPROVED

Sohil
 8/17/2017 5:59:27 PM

Quant Time: Aug 17 02:10:41 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 17 00:58:19 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2185	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	11555	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	7004	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	14794	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	15463	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	14865m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4188	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	10239	0.22	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.75	128	29996	1.067	ng/ul#	91
5) 2-Methylnaphthalene	12.36	142	22865	1.115	ng/ul	99
7) Acenaphthylene	14.24	152	11691	0.399	ng/ul	98
8) Acenaphthene	14.58	153	4806	0.209	ng/ul	96
9) Fluorene	15.57	166	8707	0.299	ng/ul	93
12) Phenanthrene	17.29	178	189508	4.704	ng/ul#	90
13) Anthracene	17.38	178	25279	0.644	ng/ul#	91
15) Fluoranthene	19.30	202	306128	6.076	ng/ul	85
17) Pyrene	19.66	202	281989	6.248	ng/ul#	81
18) Benzo(a)anthracene	21.39	228	135383	3.020	ng/ul#	85
19) Chrysene	21.44	228	149112	3.468	ng/ul	99
21) Benzo(b)fluoranthene	23.14	252	166568	3.622	ng/ul	86
22) Benzo(k)fluoranthene	23.19	252	75497m	1.665	ng/ul	
23) Benzo(a)pyrene	23.80	252	119342m	2.729	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.55	276	74066m	1.534	ng/ul	
25) Dibenzo(a,h)anthracene	26.56	278	19037m	0.472	ng/ul	
26) Benzo(a,h,i)perylene	27.37	276	61061m	1.501	ng/ul	

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

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