

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081517\
 Data File : BE093817.D
 Acq On : 15 Aug 2017 21:01
 Operator : SJ/JU
 Sample : I4672-07
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 C0AH5

Manual Integrations
 APPROVED

Sohil
 8/16/2017 6:49:57 PM

Quant Time: Aug 16 15:18:06 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 07:30:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	2427	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	12882	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	7895	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	16870	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	16670	0.40	ng/ul	0.00
20) Perylene-d12	23.91	264	12373m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4788	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	11875	0.22	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	4625	0.148	ng/ul#	94
5) 2-Methylnaphthalene	12.36	142	3229	0.141	ng/ul	97
12) Phenanthrene	17.29	178	4375	0.095	ng/ul#	91
15) Fluoranthene	19.30	202	7017	0.122	ng/ul	84
17) Pyrene	19.66	202	8479	0.174	ng/ul#	87
18) Benzo(a)anthracene	21.39	228	3018	0.062	ng/ul#	88
19) Chrysene	21.44	228	4298	0.093	ng/ul	94
21) Benzo(b)fluoranthene	23.14	252	5344m	0.140	ng/ul	
22) Benzo(k)fluoranthene	23.19	252	2163m	0.057	ng/ul	
23) Benzo(a)pyrene	23.80	252	3393m	0.093	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.56	276	1837m	0.046	ng/ul	
26) Benzo(g,h,i)perylene	27.38	276	1573m	0.046	ng/ul	

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

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