

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060217\
 Data File : BM010404.D
 Acq On : 03 Jun 2017 01:12
 Operator : SJ/MA
 Sample : I3163-22
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHSH2

Manual Integrations
 APPROVED

mohammad
 6/5/2017 11:39:11 AM

Quant Time: Jun 05 11:16:30 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM053017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 04:03:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	857	0.40	ng/ul	0.00
2) Naphthalene-d8	10.72	136	2551	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	1489	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	3189	0.40	ng/ul	0.00
16) Chrysene-d12	21.46	240	3293	0.40	ng/ul	0.00
20) Perylene-d12	23.81	264	3283m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	611	0.16	ng/ul	0.00
14) Fluoranthene-d10	19.31	212	1502	0.18	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.77	128	724	0.108	ng/ul#	81
5) 2-Methylnaphthalene	12.37	142	282	0.059	ng/ul	97
8) Acenaphthene	14.61	153	289	0.051	ng/ul#	91
9) Fluorene	15.60	166	339	0.050	ng/ul#	78
11) Pentachlorophenol	16.94	266	103	0.084	ng/ul#	85
12) Phenanthrene	17.33	178	6414	0.626	ng/ul	97
15) Fluoranthene	19.35	202	7248	0.570	ng/ul	91
17) Pyrene	19.71	202	7232	0.516	ng/ul#	95
18) Benzo(a)anthracene	21.44	228	3834	0.286	ng/ul#	87
19) Chrysene	21.49	228	6088	0.470	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.19	276	2257m	0.126	ng/ul	
26) Benzo(g,h,i)perylene	26.94	276	2820m	0.183	ng/ul	

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

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