

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025436.D
 Acq On : 10 Mar 2020 05:06
 Operator : CG/JU
 Sample : L1615-17
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBDP1

Manual Integrations
 APPROVED

mohammad
 3/11/2020 8:00:46 AM

Quant Time: Mar 10 11:45:59 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 09 13:32:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	3488	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	14779	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9888	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	21208	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	15739	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	15282	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	6245	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	14975	0.26	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	2207	0.050	ng/ul#	92
5) 2-Methylnaphthalene	12.10	142	1108	0.035	ng/ul	99
7) Acenaphthylene	14.01	152	14172	0.351	ng/ul	98
8) Acenaphthene	14.35	153	1309	0.037	ng/ul	97
9) Fluorene	15.34	166	5952	0.138	ng/ul	97
12) Phenanthrene	17.07	178	17919	0.267	ng/ul	98
13) Anthracene	17.16	178	449116	7.679	ng/ul	99
15) Fluoranthene	19.09	202	70370	0.966	ng/ul	99
17) Pyrene	19.45	202	103947	1.471	ng/ul	99
18) Benzo(a)anthracene	21.20	228	67106	1.193	ng/ul	94
19) Chrysene	21.23	228	118236	1.871	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	330294m	5.458	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	104997m	1.656	ng/ul	
23) Benzo(a)pyrene	23.31	252	144636m	2.878	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.57	276	100315	1.471	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.58	278	27585	0.481	ng/ul	97
26) Benzo(a,h,i)perylene	26.24	276	56684	0.958	ng/ul	98

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

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