

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025277.D
 Acq On : 05 Mar 2020 07:59
 Operator : CG/JU
 Sample : L1543-07
 Misc :
 ALS Vial : 73 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD71

Manual Integrations
 APPROVED

mohammad
 3/5/2020 3:31:01 PM

Quant Time: Mar 05 14:03:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 04 08:02:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3852	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	17192	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	13054	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	31548m	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	30305	0.40	ng/ul	0.00
20) Perylene-d12	23.42	264	27846	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	10677	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	33451	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.49	128	2424	0.047	ng/ul#	90
5) 2-Methylnaphthalene	12.11	142	1389	0.037	ng/ul	97
8) Acenaphthene	14.36	153	3526	0.074	ng/ul	99
9) Fluorene	15.35	166	7160	0.120	ng/ul	98
11) Pentachlorophenol	16.70	266	271	0.033	ng/ul	99
12) Phenanthrene	17.08	178	82128	0.807	ng/ul	95
13) Anthracene	17.17	178	6218	0.069	ng/ul#	95
15) Fluoranthene	19.10	202	91710	0.720	ng/ul	95
17) Pyrene	19.46	202	58519m	0.509	ng/ul	
18) Benzo(a)anthracene	21.21	228	13874	0.125	ng/ul	100
19) Chrysene	21.27	228	13973	0.114	ng/ul	98
21) Benzo(b)fluoranthene	22.77	252	11683m	0.105	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	5170m	0.044	ng/ul	
23) Benzo(a)pyrene	23.33	252	6014	0.064	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.60	276	3879	0.033	ng/ul#	100
26) Benzo(a,h,i)perylene	26.26	276	3283	0.033	ng/ul#	91

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

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