

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025153.D
 Acq On : 29 Feb 2020 03:45
 Operator : CG/JU
 Sample : L1507-21
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD19

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:39:25 AM

Quant Time: Mar 02 00:53:08 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 02 00:36:45 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2930	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	11443	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	7439	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16774m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	17253	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	18913	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	6209	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	16840	0.29	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.03	152	1387	0.032	ng/ul#	88
8) Acenaphthene	14.37	153	1833	0.064	ng/ul	98
9) Fluorene	15.36	166	1884	0.052	ng/ul	98
11) Pentachlorophenol	16.71	266	343	0.067	ng/ul	97
12) Phenanthrene	17.09	178	9000	0.153	ng/ul	99
13) Anthracene	17.19	178	2823	0.049	ng/ul#	94
15) Fluoranthene	19.11	202	22582	0.263	ng/ul	96
17) Pyrene	19.48	202	21984	0.283	ng/ul	96
18) Benzo(a)anthracene	21.22	228	7253	0.097	ng/ul	97
19) Chrysene	21.27	228	8490	0.114	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	12364m	0.167	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	4977m	0.068	ng/ul	
23) Benzo(a)pyrene	23.34	252	4307	0.064	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.62	276	3334	0.039	ng/ul#	100
26) Benzo(g,h,i)perylene	26.28	276	2506	0.035	ng/ul#	88

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

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