

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025144.D
 Acq On : 28 Feb 2020 21:45
 Operator : CG/JU
 Sample : L1507-10
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD08

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:38:43 AM

Quant Time: Feb 29 05:44:15 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 : Sat Feb 29 04:43:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	3029	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	11566	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	7186	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16743	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	17206	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	19241	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	847	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	2337	0.04	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.10	178	1954	0.033	ng/ul#	92
15) Fluoranthene	19.11	202	5775	0.068	ng/ul	99
17) Pyrene	19.47	202	6129	0.079	ng/ul	99
18) Benzo(a)anthracene	21.22	228	2735	0.037	ng/ul	95
19) Chrysene	21.27	228	3341	0.045	ng/ul	97
21) Benzo(b)fluoranthene	22.78	252	4431m	0.059	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	1822m	0.024	ng/ul	
23) Benzo(a)pyrene	23.34	252	2707	0.039	ng/ul#	73
24) Indeno(1,2,3-cd)pyrene	25.62	276	1940	0.022	ng/ul#	100
26) Benzo(g,h,i)perylene	26.28	276	1595	0.022	ng/ul#	85

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

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