

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025160.D
 Acq On : 29 Feb 2020 10:19
 Operator : CG/JU
 Sample : L1542-02
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD38

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:39:33 AM

Quant Time: Mar 02 01:37:04 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 01:15:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	2638	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	10420	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	6924	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15701m	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	17048	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	19271	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	7946	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	21488	0.40	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	1941	0.035	ng/ul#	93
15) Fluoranthene	19.11	202	4222	0.053	ng/ul	97
17) Pyrene	19.47	202	6370	0.083	ng/ul#	94
18) Benzo(a)anthracene	21.22	228	2496	0.034	ng/ul	93
19) Chrysene	21.27	228	3692	0.050	ng/ul	95
21) Benzo(b)fluoranthene	22.78	252	6051m	0.080	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	2359m	0.032	ng/ul	
23) Benzo(a)pyrene	23.34	252	2457	0.036	ng/ul#	65
24) Indeno(1,2,3-cd)pyrene	25.62	276	1900	0.022	ng/ul#	100
26) Benzo(g,h,i)perylene	26.28	276	1630	0.022	ng/ul#	76

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

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