

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022820\
 Data File : BM025128.D
 Acq On : 28 Feb 2020 10:16
 Operator : CG/JU
 Sample : L1506-22DL 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBD00DL

Manual Integrations
 APPROVED

Sohil
 3/3/2020 8:38:29 AM

Quant Time: Feb 28 11:02:29 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 : Tue Feb 25 11:03:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2994	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.46	136	11057	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.32	164	6716	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	15340	0.40	ng/ul	-0.01
16) Chrysene-d12	21.24	240	16048	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	17490	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	2342	0.11	ng/ul	-0.01
14) Fluoranthene-d10	19.09	212	6114	0.12	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.03	152	1022	0.026	ng/ul#	86
12) Phenanthrene	17.10	178	1797m	0.033	ng/ul	
13) Anthracene	17.19	178	3255	0.062	ng/ul	99
15) Fluoranthene	19.11	202	27363	0.349	ng/ul	95
17) Pyrene	19.48	202	64599	0.893	ng/ul	95
18) Benzo(a)anthracene	21.22	228	32043	0.462	ng/ul	99
19) Chrysene	21.28	228	41765	0.605	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	49082m	0.716	ng/ul	
22) Benzo(k)fluoranthene	22.83	252	23335m	0.344	ng/ul	
23) Benzo(a)pyrene	23.34	252	20679m	0.330	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.63	276	13253	0.169	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.64	278	3985	0.060	ng/ul#	68
26) Benzo(g,h,i)perylene	26.29	276	8063	0.122	ng/ul#	94

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

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