

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN072920\
 Data File : BN011292.D
 Acq On : 29 Jul 2020 14:44
 Operator : CG/JU
 Sample : L3388-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD2

Manual Integrations
 APPROVED

mohammad
 7/30/2020 3:35:31 PM

Quant Time: Jul 29 15:24:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN072020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 29 11:20:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	2133	0.40	ng/ul	0.00
2) Naphthalene-d8	10.22	136	7899	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	4583	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	10333	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	9049	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8573	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	3324	0.17	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	8596	0.19	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.90	178	753	0.022	ng/ul#	81
15) Fluoranthene	18.94	202	2965	0.078	ng/ul	93
17) Pyrene	19.30	202	2869	0.059	ng/ul#	90
18) Benzo(a)anthracene	21.07	228	1173	0.037	ng/ul#	87
19) Chrysene	21.12	228	2462	0.058	ng/ul#	94
21) Benzo(b)fluoranthene	22.59	252	3499m	0.095	ng/ul	
22) Benzo(k)fluoranthene	22.63	252	1394m	0.033	ng/ul	
23) Benzo(a)pyrene	23.12	252	1165	0.037	ng/ul#	44
24) Indeno(1,2,3-cd)pyrene	25.30	276	1170	0.032	ng/ul#	58

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

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