

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN080620\
 Data File : BN011382.D
 Acq On : 06 Aug 2020 16:50
 Operator : CG/JU
 Sample : L3487-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F4L05

Manual Integrations
 APPROVED

mohammad
 8/7/2020 1:53:15 PM

Quant Time: Aug 06 17:22:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 06 12:09:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	858	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	2834	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	1610	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	3848m	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	3472	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	4059	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.80	152	1786	0.25	ng/ul	0.00
14) Fluoranthene-d10	18.88	212	5472	0.29	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.25	128	814	0.097	ng/ul#	79
8) Acenaphthene	14.15	153	135	0.022	ng/ul#	88
9) Fluorene	15.15	166	423	0.059	ng/ul#	95
11) Pentachlorophenol	16.50	266	2847	3.018	ng/ul	100
12) Phenanthrene	16.88	178	270	0.021	ng/ul#	54
24) Indeno(1,2,3-cd)pyrene	25.25	276	321	0.021	ng/ul#	70
26) Benzo(a,h,i)perylene	25.87	276	306	0.022	ng/ul#	16

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

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