

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110719\
 Data File : BN008500.D
 Acq On : 08 Nov 2019 09:53
 Operator : JU
 Sample : K5699-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F2J00

Manual Integrations
 APPROVED

mohammad
 11/11/2019 8:08:41 AM

Quant Time: Nov 08 10:42:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 08 06:44:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	1448	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	6270	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.19	164	3612	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	7763m	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	6461	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	6962	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	2590	0.26	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	7779	0.34	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.36	128	35465	1.987	ng/ul	97
5) 2-Methylnaphthalene	11.98	142	2860	0.241	ng/ul	99
7) Acenaphthylene	13.90	152	2025	0.120	ng/ul#	93
8) Acenaphthene	14.25	153	3240	0.238	ng/ul	99
9) Fluorene	15.25	166	2071	0.139	ng/ul	97
12) Phenanthrene	16.98	178	2941	0.125	ng/ul	95
13) Anthracene	17.07	178	6525	0.316	ng/ul	99
15) Fluoranthene	19.01	202	11858	0.438	ng/ul	100
17) Pyrene	19.37	202	15207m	0.482	ng/ul	
18) Benzo(a)anthracene	21.13	228	6099	0.271	ng/ul	96
19) Chrysene	21.19	228	5449	0.190	ng/ul	97
21) Benzo(b)fluoranthene	22.67	252	14286m	0.551	ng/ul	
22) Benzo(k)fluoranthene	22.71	252	4383m	0.149	ng/ul	
23) Benzo(a)pyrene	23.22	252	5523	0.211	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.46	276	3536	0.119	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.46	278	966	0.041	ng/ul#	52
26) Benzo(a,h,i)perylene	26.10	276	3000	0.113	ng/ul	93

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

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