

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122819\
 Data File : BN009224.D
 Acq On : 28 Dec 2019 12:42
 Operator : JU
 Sample : K6278-06DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0C53DL

Manual Integrations
 APPROVED

mohammad
 12/29/2019 5:32:13 PM

Quant Time: Dec 29 04:18:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Dec 29 03:34:38 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	2045	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	7401	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.14	164	4027	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.89	188	8091	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	6452	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	6594	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.86	152	322	0.02	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	699	0.03	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.31	128	2969	0.132	ng/ul#	94
5) 2-Methylnaphthalene	11.94	142	1843	0.109	ng/ul	100
7) Acenaphthylene	13.86	152	1682	0.080	ng/ul#	91
9) Fluorene	15.20	166	487	0.025	ng/ul#	93
12) Phenanthrene	16.93	178	11599	0.412	ng/ul	99
13) Anthracene	17.02	178	2342	0.093	ng/ul#	93
15) Fluoranthene	18.95	202	27611	0.858	ng/ul	99
17) Pyrene	19.32	202	22727	0.690	ng/ul	99
18) Benzo(a)anthracene	21.07	228	12135	0.424	ng/ul	96
19) Chrysene	21.12	228	14060	0.488	ng/ul	96
21) Benzo(b)fluoranthene	22.58	252	18372m	0.597	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	7222m	0.233	ng/ul	
23) Benzo(a)pyrene	23.12	252	10799	0.400	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.30	276	8224	0.280	ng/ul	90
25) Dibenzo(a,h)anthracene	25.31	278	2193	0.094	ng/ul#	65
26) Benzo(a,h,i)perylene	25.93	276	6478	0.262	ng/ul	95

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

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