

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120219\
 Data File : BN008815.D
 Acq On : 02 Dec 2019 15:30
 Operator : JU
 Sample : K5851-01DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AA1DL

Manual Integrations
 APPROVED

mohammad
 12/3/2019 2:38:52 PM

Quant Time: Dec 02 16:28:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 02 12:46:36 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	5364	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	22266	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	12384	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	25184	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	19054	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	16683	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	1644	0.05	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	3704	0.05	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.39	128	3121	0.046	ng/ul#	94
5) 2-Methylnaphthalene	12.02	142	1777	0.038	ng/ul	98
7) Acenaphthylene	13.93	152	6765	0.099	ng/ul	98
9) Fluorene	15.27	166	1403	0.024	ng/ul#	93
12) Phenanthrene	17.00	178	9841	0.116	ng/ul	99
13) Anthracene	17.09	178	4539	0.057	ng/ul#	94
15) Fluoranthene	19.02	202	40415	0.387	ng/ul	98
17) Pyrene	19.39	202	35405	0.463	ng/ul	98
18) Benzo(a)anthracene	21.15	228	23513	0.307	ng/ul#	89
19) Chrysene	21.20	228	22402m	0.298	ng/ul	
21) Benzo(b)fluoranthene	22.68	252	27133m	0.394	ng/ul	
22) Benzo(k)fluoranthene	22.72	252	9093m	0.133	ng/ul	
23) Benzo(a)pyrene	23.23	252	18441	0.286	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.45	276	11179	0.148	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.46	278	2680	0.044	ng/ul#	1
26) Benzo(a,h,i)perylene	26.09	276	11329	0.179	ng/ul#	84

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

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