

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102419\
 Data File : BN008412.D
 Acq On : 25 Oct 2019 05:44
 Operator : JU
 Sample : K5236-06DL 5X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPL6DL

Manual Integrations
 APPROVED

mohammad
 10/25/2019 9:28:11 PM

Quant Time: Oct 25 12:01:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 25 00:45:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1582	0.40	ng/ul	0.00
2) Naphthalene-d8	10.34	136	8027	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.21	164	5450	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	11977	0.40	ng/ul	0.00
16) Chrysene-d12	21.17	240	11325	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	9427	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	410	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	1457	0.04	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.93	152	1783	0.069	ng/ul#	91
12) Phenanthrene	17.00	178	8722	0.249	ng/ul	99
13) Anthracene	17.09	178	1957	0.063	ng/ul#	86
15) Fluoranthene	19.03	202	18121	0.391	ng/ul	98
17) Pyrene	19.39	202	14417	0.320	ng/ul	98
18) Benzo(a)anthracene	21.15	228	8705	0.211	ng/ul	91
19) Chrysene	21.20	228	8195m	0.162	ng/ul	
21) Benzo(b)fluoranthene	22.69	252	10449m	0.340	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	4156m	0.120	ng/ul	
23) Benzo(a)pyrene	23.24	252	6328	0.196	ng/ul#	72
24) Indeno(1,2,3-cd)pyrene	25.49	276	4721	0.124	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.49	278	1013	0.034	ng/ul#	3
26) Benzo(g,h,i)perylene	26.13	276	4668	0.134	ng/ul#	84

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

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