

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062419\
 Data File : BN006561.D
 Acq On : 25 Jun 2019 08:43
 Operator : HP/JU
 Sample : K3360-14DL 5X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41T6DL

Manual Integrations
 APPROVED

mohammad
 6/25/2019 10:21:41 PM

Quant Time: Jun 25 11:48:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 25 07:15:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	5458	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	19874	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	9477	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	17265	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	10354	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	13679	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	946	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	1585	0.03	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.46	128	2638	0.048	ng/ul#	91
5) 2-Methylnaphthalene	12.08	142	1305	0.034	ng/ul	100
8) Acenaphthene	14.34	153	13305	0.364	ng/ul	99
9) Fluorene	15.34	166	18795	0.440	ng/ul	98
12) Phenanthrene	17.08	178	317589	5.730	ng/ul	99
13) Anthracene	17.17	178	37398	0.707	ng/ul	99
15) Fluoranthene	19.11	202	404098	6.603	ng/ul	100
17) Pyrene	19.48	202	296253	5.432	ng/ul	97
18) Benzo(a)anthracene	21.24	228	119273	2.605	ng/ul	99
19) Chrysene	21.29	228	134228	3.122	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	149352m	2.618	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	61101m	1.100	ng/ul	
23) Benzo(a)pyrene	23.39	252	100033	1.883	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.72	276	67180	1.168	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.72	278	17089	0.372	ng/ul	99
26) Benzo(g,h,i)perylene	26.41	276	58578	1.216	ng/ul	95

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

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