

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062419\
 Data File : BN006567.D
 Acq On : 25 Jun 2019 12:40
 Operator : HP/JU
 Sample : K3360-21MS
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4202MS

Manual Integrations
 APPROVED

mohammad
 6/25/2019 10:21:58 PM

Quant Time: Jun 25 13:22:33 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	4027	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	14094	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	6385	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	11614	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	8857	0.40	ng/ul	0.03
20) Perylene-d12	23.54	264	10392	0.40	ng/ul	0.06
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	4907	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	8485	0.27	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.46	128	914	0.024	ng/ul#	74
7) Acenaphthylene	14.01	152	979	0.028	ng/ul#	78
8) Acenaphthene	14.35	153	6992	0.284	ng/ul	99
9) Fluorene	15.35	166	816	0.028	ng/ul#	93
11) Pentachlorophenol	16.74	266	110	0.037	ng/ul	92
12) Phenanthrene	17.09	178	30661	0.822	ng/ul	99
13) Anthracene	17.18	178	4932	0.139	ng/ul#	94
15) Fluoranthene	19.12	202	130810	3.178	ng/ul	100
17) Pyrene	19.48	202	120129	2.575	ng/ul	100
18) Benzo(a)anthracene	21.27	228	54372	1.388	ng/ul	100
19) Chrysene	21.32	228	73604	2.001	ng/ul	100
21) Benzo(b)fluoranthene	22.87	252	109137m	2.519	ng/ul	
22) Benzo(k)fluoranthene	22.91	252	40791m	0.966	ng/ul	
23) Benzo(a)pyrene	23.44	252	63314	1.569	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.78	276	51911	1.188	ng/ul#	87
25) Dibenzo(a,h)anthracene	25.79	278	11820	0.338	ng/ul	99
26) Benzo(g,h,i)perylene	26.47	276	44775	1.224	ng/ul	94

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

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