

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061719\
 Data File : BN006344.D
 Acq On : 17 Jun 2019 12:58
 Operator : JU/SJ
 Sample : K3332-16MSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4222MSD

Manual Integrations
 APPROVED

mohammad
 6/20/2019 10:50:51 AM

Quant Time: Jun 18 02:00:11 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 : Mon Jun 17 11:13:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	3815	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15242	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7971	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	14005	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	9497	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	10649	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	4739	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	9037	0.24	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.36	153	7392	0.241	ng/ul	99
11) Pentachlorophenol	16.72	266	301	0.085	ng/ul	89
12) Phenanthrene	17.09	178	2749	0.061	ng/ul#	91
15) Fluoranthene	19.12	202	9375	0.189	ng/ul	94
17) Pyrene	19.49	202	21572	0.431	ng/ul	99
18) Benzo(a)anthracene	21.26	228	3992	0.095	ng/ul	93
19) Chrysene	21.31	228	4004	0.102	ng/ul	95
21) Benzo(b)fluoranthene	22.84	252	6266m	0.141	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	2470m	0.057	ng/ul	
23) Benzo(a)pyrene	23.41	252	3604	0.087	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	25.75	276	2908	0.065	ng/ul#	88
26) Benzo(g,h,i)perylene	26.43	276	2396	0.064	ng/ul#	82

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

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