

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121119\
 Data File : BP001324.D
 Acq On : 11 Dec 2019 21:27
 Operator : JU
 Sample : K6204-06 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 K072-AA-WC-1

Manual Integrations
 APPROVED

mohammad
 12/12/2019 11:34:26 AM

Quant Time: Dec 12 08:07:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 05 12:34:47 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	55295	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	194805	20.00	ng	-0.01
39) Acenaphthene-d10	13.20	164	114110	20.00	ng	0.00
64) Phenanthrene-d10	15.60	188	214966	20.00	ng	-0.01
76) Chrysene-d12	19.24	240	171259	20.00	ng	-0.01
87) Perylene-d12	21.02	264	191204	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	118020	35.41	ng	0.00
7) Phenol-d6	7.75	99	185933	41.87	ng	-0.01
23) Nitrobenzene-d5	9.18	82	124061	27.63	ng	-0.02
42) 2,4,6-Tribromophenol	14.49	330	80512	73.61	ng	-0.01
45) 2-Fluorobiphenyl	12.12	172	287478	36.87	ng	-0.02
79) Terphenyl-d14	17.88	244	408021	44.08	ng	-0.01
Target Compounds						Qvalue
49) Acenaphthylene	12.96	152	21338	2.021	ng	98
52) Acenaphthene	13.25	154	20370	3.207	ng	94
58) Fluorene	14.10	166	27857	3.642	ng	96
71) Phenanthrene	15.64	178	499522	44.200	ng	99
72) Anthracene	15.71	178	115900	10.405	ng	99
73) Carbazole	15.96	167	50029	4.936	ng	98
75) Fluoranthene	17.35	202	785003	65.612	ng	99
78) Pyrene	17.66	202	638177	47.312	ng	98
80) Butylbenzylphthalate	18.53	149	13072	2.098	ng	95
81) Benzo(a)anthracene	19.23	228	336882	28.371	ng	98
83) Chrysene	19.27	228	306933	28.867	ng	97
86) Indeno(1,2,3-cd)pyrene	22.66	276	185424m	14.797	ng	
88) Benzo(b)fluoranthene	20.53	252	394679m	33.795	ng	
89) Benzo(k)fluoranthene	20.56	252	135042m	12.006	ng	
90) Benzo(a)pyrene	20.94	252	271415	25.231	ng	96
91) Dibenzo(a,h)anthracene	22.67	278	47989m	4.559	ng	
92) Benzo(a,h,i)perylene	23.14	276	163165	15.696	ng	97

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

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