

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082319\
 Data File : BF116478.D
 Acq On : 23 Aug 2019 15:12
 Operator : HP/JU
 Sample : K4418-01RE
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CFVP-228LRE

Manual Integrations
 APPROVED

Jagrut
 8/28/2019 8:09:33 AM

Quant Time: Aug 23 17:04:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 23 13:38:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	111946	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	445596	20.00	ng	0.00
39) Acenaphthene-d10	9.92	164	214063	20.00	ng	0.00
64) Phenanthrene-d10	11.40	188	268950	20.00	ng	0.00
76) Chrysene-d12	14.03	240	192136	20.00	ng	0.00
87) Perylene-d12	15.50	264	190006	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	93645	12.22	ng	0.00
7) Phenol-d6	6.49	99	81452	8.67	ng	-0.01
23) Nitrobenzene-d5	7.43	82	314234	38.58	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	58395	31.77	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	584192	43.95	ng	0.00
79) Terphenyl-d14	12.98	244	401906	39.07	ng	0.00
Target Compounds						
10) Phenol	6.51	94	37999	3.725	ng	96
20) 3+4-Methylphenols	7.27	107	84552	10.361	ng	# 69
32) Benzoic acid	7.87	122	16124m	3.923	ng	
50) Dimethylphthalate	9.63	163	232165	15.545	ng	99
71) Phenanthrene	11.42	178	64686	4.415	ng	99
75) Fluoranthene	12.60	202	55598	3.670	ng	95
78) Pyrene	12.83	202	69157	4.501	ng	98
81) Benzo(a)anthracene	14.02	228	31072	2.497	ng	# 80
83) Chrysene	14.06	228	30142	2.390	ng	# 94
88) Benzo(b)fluoranthene	15.07	252	41674m	3.550	ng	
90) Benzo(a)pyrene	15.43	252	38293	3.691	ng	# 87
92) Benzo(a,h,i)perylene	17.39	276	21874	2.692	ng	# 86

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

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