

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090119\
 Data File : BF116750.D
 Acq On : 1 Sep 2019 18:42
 Operator : HP/JU
 Sample : K4547-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T3-S-B1-B4

Manual Integrations
 APPROVED

mohammad
 9/5/2019 9:54:41 AM

Quant Time: Sep 02 06:04:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	84098	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	333771	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	180287	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	333907	20.00	ng	-0.01
76) Chrysene-d12	13.99	240	253764	20.00	ng	0.00
87) Perylene-d12	15.44	264	258063	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	424390	81.75	ng	0.00
7) Phenol-d6	6.47	99	589995	86.55	ng	-0.01
23) Nitrobenzene-d5	7.40	82	346884	59.33	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	117148	97.20	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	576735	53.96	ng	-0.01
79) Terphenyl-d14	12.94	244	690795	50.36	ng	-0.01
Target Compounds						
10) Phenol	6.48	94	44198	5.856	ng	95
50) Dimethylphthalate	9.59	163	90781	7.293	ng	97
71) Phenanthrene	11.38	178	43263	2.328	ng	99
75) Fluoranthene	12.57	202	122624	6.713	ng	98
78) Pyrene	12.80	202	118569	5.256	ng	98
81) Benzo(a)anthracene	13.98	228	58550m	3.646	ng	
83) Chrysene	14.02	228	65026	3.930	ng	96
84) Bis(2-ethylhexyl)phthalate	13.97	149	110650	8.171	ng	97
86) Indeno(1,2,3-cd)pyrene	16.88	276	33755	2.540	ng	95
88) Benzo(b)fluoranthene	15.02	252	87104m	5.583	ng	
90) Benzo(a)pyrene	15.37	252	53741	3.959	ng	# 94
92) Benzo(a,h,i)perylene	17.32	276	29373	2.424	ng	# 94

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

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