

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103119\
 Data File : BF117671.D
 Acq On : 31 Oct 2019 21:07
 Operator : JU
 Sample : K5148-13 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-102919

Manual Integrations
 APPROVED

mohammad
 11/1/2019 12:34:55 PM

Quant Time: Nov 01 02:21:11 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	
1) 1,4-Dichlorobenzene-d4	6.73	152	37448	20.00	ng	0.00	
21) Naphthalene-d8	8.01	136	164253	20.00	ng	0.00	
39) Acenaphthene-d10	9.76	164	76215	20.00	ng	0.00	
64) Phenanthrene-d10	11.25	188	102589	20.00	ng	0.00	
76) Chrysene-d12	13.90	240	77474	20.00	ng	0.00	
87) Perylene-d12	15.34	264	66987	20.00	ng	0.02	
System Monitoring Compounds							
5) 2-Fluorophenol	5.38	112	36996	14.70	ng	0.03	
7) Phenol-d6	6.41	99	59914	17.72	ng	0.03	
23) Nitrobenzene-d5	7.32	82	32714	9.83	ng	0.01	
42) 2,4,6-Tribromophenol	10.57	330	8293	12.58	ng	0.00	
45) 2-Fluorobiphenyl	9.09	172	60436	11.17	ng	0.00	
79) Terphenyl-d14	12.83	244	42352	8.92	ng	0.00	
Target Compounds							
71) Phenanthrene	11.27	178	34963	5.869	ng		99
72) Anthracene	11.34	178	14019m	2.294	ng		
75) Fluoranthene	12.47	202	73043	12.249	ng		99
78) Pyrene	12.70	202	68444	10.285	ng		99
81) Benzo(a)anthracene	13.89	228	26265	5.026	ng		98
83) Chrysene	13.93	228	25719	5.022	ng		96
86) Indeno(1,2,3-cd)pyrene	16.78	276	7850	2.053	ng		96
88) Benzo(b)fluoranthene	14.93	252	20363m	5.148	ng		
89) Benzo(k)fluoranthene	14.96	252	9800m	2.417	ng		
90) Benzo(a)pyrene	15.29	252	16875	4.692	ng	#	91
92) Benzo(a,h,i)perylene	17.21	276	7375	2.506	ng		95

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

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