

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\  
 Data File : BF120192.D  
 Acq On : 24 Apr 2020 10:39  
 Operator : MA/SJ  
 Sample : L2397-05  
 Misc :  
 ALS Vial : 29 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-1-COMP

Manual Integrations  
 APPROVED

mohammad  
 4/27/2020 3:04:07 PM

Quant Time: Apr 24 11:22:15 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Fri Apr 24 10:24:30 2020  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	
1) 1,4-Dichlorobenzene-d4	6.66	152	141062	20.00	ng	0.00	
21) Naphthalene-d8	7.95	136	552798	20.00	ng	0.00	
39) Acenaphthene-d10	9.70	164	284516	20.00	ng	0.00	
64) Phenanthrene-d10	11.18	188	533545	20.00	ng	0.00	
76) Chrysene-d12	13.82	240	336165	20.00	ng	-0.01	
87) Perylene-d12	15.22	264	283821	20.00	ng	0.00	
System Monitoring Compounds							
5) 2-Fluorophenol	5.27	112	832930	102.04	ng	0.00	
7) Phenol-d6	6.31	99	1107233	105.27	ng	0.00	
23) Nitrobenzene-d5	7.23	82	724523	67.80	ng	-0.01	
42) 2,4,6-Tribromophenol	10.49	330	311494	89.42	ng	0.00	
45) 2-Fluorobiphenyl	9.03	172	1450612	72.39	ng	0.00	
79) Terphenyl-d14	12.77	244	1543039	77.24	ng	0.00	
Target Compounds							
50) Dimethylphthalate	9.42	163	201801	9.170	ng		100
71) Phenanthrene	11.20	178	75122	2.646	ng		97
75) Fluoranthene	12.39	202	207233	6.764	ng		94
78) Pyrene	12.62	202	249013	8.787	ng		98
81) Benzo(a)anthracene	13.80	228	116167	5.253	ng		94
83) Chrysene	13.84	228	106241	4.728	ng		97
86) Indeno(1,2,3-cd)pyrene	16.56	276	53321	2.692	ng		96
88) Benzo(b)fluoranthene	14.82	252	126601m	6.201	ng		
89) Benzo(k)fluoranthene	14.84	252	38005m	2.157	ng		
90) Benzo(a)pyrene	15.16	252	98214	5.721	ng		99
92) Benzo(a,h,i)perylene	16.96	276	48156	3.208	ng	#	94

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

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