

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060519\  
 Data File : BF114828.D  
 Acq On : 5 Jun 2019 19:54  
 Operator : HP/JU  
 Sample : K2638-02  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 2019

Manual Integrations  
 APPROVED

mohammad  
 6/6/2019 4:50:12 PM

Quant Time: Jun 06 06:55:26 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF060419.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Jun 04 16:57:15 2019  
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	
1) 1,4-Dichlorobenzene-d4	6.67	152	345561	20.00	ng	0.00	
21) Naphthalene-d8	7.95	136	1358189	20.00	ng	0.00	
39) Acenaphthene-d10	9.70	164	701151	20.00	ng	0.00	
64) Phenanthrene-d10	11.17	188	1388445	20.00	ng	0.00	
76) Chrysene-d12	13.79	240	963907	20.00	ng	0.00	
87) Perylene-d12	15.19	264	1012525	20.00	ng	0.01	
System Monitoring Compounds							
5) 2-Fluorophenol	5.27	112	1963028	99.26	ng	0.01	
7) Phenol-d6	6.31	99	2436791	101.26	ng	0.00	
23) Nitrobenzene-d5	7.23	82	1513916	68.62	ng	0.00	
42) 2,4,6-Tribromophenol	10.48	330	770698	102.00	ng	0.00	
45) 2-Fluorobiphenyl	9.03	172	2698264	72.09	ng	0.00	
79) Terphenyl-d14	12.76	244	3080452	62.15	ng	0.00	
Target Compounds							
50) Dimethylphthalate	9.42	163	660329	12.895	ng		99
71) Phenanthrene	11.19	178	281231	4.243	ng		95
75) Fluoranthene	12.37	202	609115	8.829	ng		96
78) Pyrene	12.60	202	548005	6.554	ng		98
81) Benzo(a)anthracene	13.79	228	290450	3.910	ng		98
83) Chrysene	13.82	228	261445	3.618	ng		95
88) Benzo(b)fluoranthene	14.80	252	366707m	5.689	ng		
90) Benzo(a)pyrene	15.13	252	258321	4.630	ng	#	96
92) Benzo(g,h,i)perylene	16.88	276	137785	2.593	ng		97

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

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