

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
 Data File : BF118388.D
 Acq On : 30 Dec 2019 18:37
 Operator : JU
 Sample : K6441-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K321-WC-02

Manual Integrations
 APPROVED

mohammad
 12/31/2019 12:04:15 PM

mohammad

Quant Time: Dec 31 01:30:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	12/31/2019 12:04:15 PM
1) 1,4-Dichlorobenzene-d4	6.85	152	131013	20.00	ng	0.00	
21) Naphthalene-d8	8.14	136	528047	20.00	ng	0.00	
39) Acenaphthene-d10	9.90	164	288630	20.00	ng	0.00	
64) Phenanthrene-d10	11.39	188	535899	20.00	ng	0.00	
76) Chrysene-d12	14.05	240	323643	20.00	ng	0.00	
87) Perylene-d12	15.54	264	339997	20.00	ng	0.00	
System Monitoring Compounds							
5) 2-Fluorophenol	5.47	112	701465	91.39	ng	0.00	
7) Phenol-d6	6.49	99	891911	93.68	ng	0.00	
23) Nitrobenzene-d5	7.42	82	539055	58.39	ng	0.00	
42) 2,4,6-Tribromophenol	10.70	330	319604	89.67	ng	0.00	
45) 2-Fluorobiphenyl	9.22	172	1135997	60.08	ng	0.00	
79) Terphenyl-d14	12.99	244	1313073	67.31	ng	0.00	
Target Compounds							
50) Dimethylphthalate	9.62	163	97465	4.499	ng	100	
71) Phenanthrene	11.42	178	749768	25.664	ng	99	
72) Anthracene	11.47	178	195554	6.687	ng	98	
75) Fluoranthene	12.62	202	1018395	34.541	ng	97	
78) Pyrene	12.84	202	877353	33.197	ng	99	
81) Benzo(a)anthracene	14.04	228	391697	17.290	ng	99	
83) Chrysene	14.07	228	326625	15.058	ng	97	
86) Indeno(1,2,3-cd)pyrene	17.05	276	193301	10.459	ng	94	
88) Benzo(b)fluoranthene	15.10	252	408790m	18.113	ng		
89) Benzo(k)fluoranthene	15.13	252	140382m	6.478	ng		
90) Benzo(a)pyrene	15.48	252	296320	14.904	ng	98	
91) Dibenzo(a,h)anthracene	17.06	278	50590m	2.932	ng		
92) Benzo(a,h,i)perylene	17.51	276	180673	10.841	ng	98	

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

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