

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081719\
 Data File : BF116369.D
 Acq On : 17 Aug 2019 20:34
 Operator : HP/JU
 Sample : K4223-25
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-13

Manual Integrations
 APPROVED

Jagrut
 8/19/2019 3:31:16 PM

Quant Time: Aug 19 07:32:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	134509	20.00	ng	-0.04
21) Naphthalene-d8	8.09	136	514434	20.00	ng	-0.04
39) Acenaphthene-d10	9.84	164	255858	20.00	ng	-0.05
64) Phenanthrene-d10	11.33	188	386502	20.00	ng	-0.04
76) Chrysene-d12	13.97	240	243525	20.00	ng	-0.04
87) Perylene-d12	15.43	264	276550	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	758859	94.45	ng	-0.03
7) Phenol-d6	6.46	99	977846	96.46	ng	-0.03
23) Nitrobenzene-d5	7.37	82	614590	68.96	ng	-0.04
42) 2,4,6-Tribromophenol	10.63	330	220360	88.83	ng	-0.04
45) 2-Fluorobiphenyl	9.16	172	935724	68.50	ng	-0.04
79) Terphenyl-d14	12.90	244	754017	65.24	ng	-0.05
Target Compounds						
10) Phenol	6.47	94	24475	2.050	ng	91
50) Dimethylphthalate	9.56	163	206700	11.865	ng	99
75) Fluoranthene	12.54	202	60750	2.937	ng	97
78) Pyrene	12.77	202	49922	2.719	ng	98
88) Benzo(b)fluoranthene	15.01	252	34736m	2.172	ng	

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

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