

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010320\
 Data File : BF118472.D
 Acq On : 3 Jan 2020 22:26
 Operator : JU
 Sample : K6471-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MC-123019-2-5-COMP-B8

Manual Integrations
 APPROVED

mohammad
 1/6/2020 11:44:19 AM

Quant Time: Jan 04 01:28:43 2020
 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)
1) 1,4-Dichlorobenzene-d4	6.85	152	70090	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	255377	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	117230	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	234967	20.00	ng	0.00
76) Chrysene-d12	14.04	240	172018	20.00	ng	0.00
87) Perylene-d12	15.53	264	197314	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	314294	76.54	ng	0.00
7) Phenol-d6	6.49	99	390645	76.69	ng	0.00
23) Nitrobenzene-d5	7.41	82	243835	54.61	ng	-0.01
42) 2,4,6-Tribromophenol	10.69	330	111382	76.94	ng	0.00
45) 2-Fluorobiphenyl	9.21	172	427607	55.68	ng	-0.01
79) Terphenyl-d14	12.97	244	450425	43.44	ng	-0.01
Target Compounds						
10) Phenol	6.50	94	13184	2.303	ng	Qvalue # 76
50) Dimethylphthalate	9.61	163	37862	4.303	ng	100
75) Fluoranthene	12.61	202	42656	3.300	ng	95
78) Pyrene	12.84	202	42680	3.038	ng	98
81) Benzo(a)anthracene	14.03	228	27190	2.258	ng	98
83) Chrysene	14.07	228	25494	2.211	ng	94
88) Benzo(b)fluoranthene	15.10	252	31855m	2.432	ng	
90) Benzo(a)pyrene	15.47	252	23326	2.022	ng	# 70

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

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