

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082819\
 Data File : BF116622.D
 Acq On : 28 Aug 2019 18:39
 Operator : HP/JU
 Sample : K4470-12
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S703-N-1.0-1.5-082119

Manual Integrations
 APPROVED

Jagrut
 8/29/2019 7:46:55 PM

Quant Time: Aug 29 02:14:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 28 13:34:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	109882	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	381387	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	199999	20.00	ng	-0.01
64) Phenanthrene-d10	11.37	188	359153	20.00	ng	0.00
76) Chrysene-d12	14.00	240	277047	20.00	ng	0.00
87) Perylene-d12	15.46	264	251268	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	904012	142.10	ng	0.00
7) Phenol-d6	6.49	99	1322281	153.31	ng	0.00
23) Nitrobenzene-d5	7.42	82	788930	118.36	ng	0.00
42) 2,4,6-Tribromophenol	10.68	330	277226	145.66	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1279910	106.50	ng	0.00
79) Terphenyl-d14	12.96	244	1115375	65.48	ng	0.00
Target Compounds						
10) Phenol	6.50	94	32376	3.422	ng	99
50) Dimethylphthalate	9.60	163	222361	16.538	ng	100
71) Phenanthrene	11.40	178	42784	2.218	ng	99
75) Fluoranthene	12.58	202	94257	5.448	ng	97
78) Pyrene	12.81	202	80485	3.159	ng	98
81) Benzo(a)anthracene	13.99	228	38956	2.109	ng	99
86) Indeno(1,2,3-cd)pyrene	16.90	276	14290	4.803	ng	96
88) Benzo(b)fluoranthene	15.03	252	38330m	2.484	ng	

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

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