

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012220\
 Data File : BF118674.D
 Acq On : 22 Jan 2020 20:47
 Operator : CG/JU
 Sample : L1213-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-2

Manual Integrations
 APPROVED

mohammad
 1/23/2020 1:54:26 PM

Quant Time: Jan 23 07:21:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 20 14:45:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	335340	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	1259153	20.00	ng	-0.01
39) Acenaphthene-d10	9.92	164	672590	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	1135261	20.00	ng	0.00
76) Chrysene-d12	14.04	240	830516	20.00	ng	0.00
87) Perylene-d12	15.52	264	972733	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1755426	97.22	ng	0.00
7) Phenol-d6	6.50	99	2188406	93.25	ng	0.00
23) Nitrobenzene-d5	7.43	82	1420880	63.12	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	689041	88.57	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	2536600	61.99	ng	-0.01
79) Terphenyl-d14	12.99	244	2440195	64.08	ng	0.00
Target Compounds						
10) Phenol	6.52	94	160903	6.016	ng	90
31) Naphthalene	8.18	128	210216	3.696	ng	95
50) Dimethylphthalate	9.63	163	302092	6.894	ng	95
52) Acenaphthene	9.95	154	163809	4.457	ng	97
55) Dibenzofuran	10.12	168	142289	2.794	ng	95
58) Fluorene	10.46	166	163557	4.090	ng	98
71) Phenanthrene	11.42	178	468303	8.550	ng	93
75) Fluoranthene	12.62	202	341981	5.920	ng	94
78) Pyrene	12.84	202	267189	4.739	ng	92
83) Chrysene	14.06	228	101495	2.120	ng	94
88) Benzo(b)fluoranthene	15.08	252	123497m	2.020	ng	

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

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