

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012720\
 Data File : BF118724.D
 Acq On : 27 Jan 2020 17:37
 Operator : CG/JU
 Sample : L1010-10 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-012420

Manual Integrations
 APPROVED

mohammad
 1/28/2020 1:13:19 PM

Quant Time: Jan 27 18:12:39 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 27 14:24:10 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	287333	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	1013211	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	482682	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	727652	20.00	ng	0.00
76) Chrysene-d12	14.02	240	622899	20.00	ng	0.00
87) Perylene-d12	15.49	264	729001	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	183189	11.84	ng	-0.01
7) Phenol-d6	6.48	99	224927	11.19	ng	-0.01
23) Nitrobenzene-d5	7.41	82	124557	6.88	ng	-0.01
42) 2,4,6-Tribromophenol	10.68	330	46482	8.33	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	273987	9.33	ng	0.00
79) Terphenyl-d14	12.96	244	247927	8.68	ng	0.00
Target Compounds						
71) Phenanthrene	11.40	178	84570	2.409	ng	96
75) Fluoranthene	12.59	202	173676	4.691	ng	96
78) Pyrene	12.82	202	160366	3.792	ng	97
81) Benzo(a)anthracene	14.01	228	84160	2.246	ng	# 93
83) Chrysene	14.04	228	76804	2.139	ng	# 93
88) Benzo(b)fluoranthene	15.06	252	102236m	2.231	ng	

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

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