

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP010320\
 Data File : BP001527.D
 Acq On : 04 Jan 2020 02:16
 Operator : JU
 Sample : L1011-02 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MC-010220-0-2-COMP-B3

Quant Time: Jan 04 04:00:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 03 15:57:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	61218	20.00	ng	-0.01
21) Naphthalene-d8	10.47	136	230942	20.00	ng	-0.01
39) Acenaphthene-d10	14.33	164	150044	20.00	ng	-0.01
64) Phenanthrene-d10	17.09	188	338510	20.00	ng	-0.01
76) Chrysene-d12	21.19	240	303492	20.00	ng	-0.02
87) Perylene-d12	23.44	264	324730	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	72279	19.84	ng	0.02
7) Phenol-d6	6.90	99	101914	20.76	ng	0.02
23) Nitrobenzene-d5	8.84	82	63015	13.48	ng	0.00
42) 2,4,6-Tribromophenol	15.83	330	33653	19.82	ng	-0.01
45) 2-Fluorobiphenyl	12.95	172	131831	12.92	ng	-0.02
79) Terphenyl-d14	19.67	244	201183	13.21	ng	-0.01
Target Compounds						Qvalue
71) Phenanthrene	17.13	178	60646	3.303	ng	99
75) Fluoranthene	19.11	202	110024	4.950	ng	99
78) Pyrene	19.46	202	107351	4.874	ng	98
81) Benzo(a)anthracene	21.17	228	53029	2.490	ng	96
83) Chrysene	21.22	228	50553	2.476	ng	95
88) Benzo(b)fluoranthene	22.76	252	60487	3.008	ng	# 88
90) Benzo(a)pyrene	23.33	252	45612	2.423	ng	94

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

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