

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
 Data File : BF120459.D
 Acq On : 30 May 2020 00:23
 Operator : MA/SJ
 Sample : L2792-11 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM85-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:21:57 PM

Quant Time: May 30 05:33:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 12:39:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	242973	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	865833	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	383724	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	602677	20.00	ng	0.00
76) Chrysene-d12	13.99	240	510832	20.00	ng	0.00
86) Perylene-d12	15.44	264	491907	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	294605	23.11	ng	0.00
7) Phenol-d6	6.46	99	366739	23.34	ng	-0.01
23) Nitrobenzene-d5	7.39	82	219619	16.59	ng	-0.01
42) 2,4,6-Tribromophenol	10.66	330	74614	20.99	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	458304	20.26	ng	0.00
79) Terphenyl-d14	12.94	244	379966	16.55	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	61395	2.559	ng	98
71) Phenanthrene	11.38	178	55643	2.095	ng	96
75) Fluoranthene	12.57	202	138808	4.858	ng	96
78) Pyrene	12.80	202	131128	3.691	ng	97
81) Benzo(a)anthracene	13.99	228	81425	2.974	ng	97
83) Chrysene	14.02	228	74804	2.606	ng	97
88) Benzo(b)fluoranthene	15.02	252	92315m	3.379	ng	
90) Benzo(a)pyrene	15.38	252	63384	2.652	ng	# 96

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

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