

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060320\
 Data File : BF120519.D
 Acq On : 3 Jun 2020 19:42
 Operator : JU/CG
 Sample : L2839-17
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM108-2-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/4/2020 12:12:46 PM

Quant Time: Jun 04 01:18:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 03 14:26:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	158891	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	633262	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	289975	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	536050	20.00	ng	0.00
76) Chrysene-d12	14.00	240	404276	20.00	ng	-0.01
86) Perylene-d12	15.45	264	353491	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	853441	102.36	ng	0.00
7) Phenol-d6	6.47	99	994279	96.75	ng	0.00
23) Nitrobenzene-d5	7.40	82	718014	74.17	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	287622	107.05	ng	0.00
45) 2-Fluorobiphenyl	9.19	172	1109110	64.90	ng	0.00
79) Terphenyl-d14	12.94	244	1004107	55.27	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	178411	9.841	ng	99
71) Phenanthrene	11.38	178	58164	2.462	ng	97
75) Fluoranthene	12.57	202	149273	5.873	ng	99
78) Pyrene	12.80	202	130634	4.646	ng	97
81) Benzo(a)anthracene	13.99	228	97755	4.512	ng	95
83) Chrysene	14.02	228	93547m	4.118	ng	
88) Benzo(b)fluoranthene	15.03	252	114699m	5.842	ng	
90) Benzo(a)pyrene	15.39	252	73211	4.263	ng	98
92) Benzo(g,h,i)perylene	17.33	276	33142	2.132	ng	96

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

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