

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF052920\
 Data File : BF120452.D
 Acq On : 29 May 2020 21:12
 Operator : MA/SJ
 Sample : L2792-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NM84-COMP-2020-0-6

Manual Integrations
 APPROVED

mohammad
 6/1/2020 12:21:29 PM

Quant Time: May 30 04:39:13 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	289668	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1050656	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	497006	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	793062	20.00	ng	0.00
76) Chrysene-d12	13.99	240	564510	20.00	ng	0.00
86) Perylene-d12	15.44	264	589710	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1145762	75.38	ng	0.00
7) Phenol-d6	6.46	99	1529679	81.65	ng	0.00
23) Nitrobenzene-d5	7.40	82	999783	62.25	ng	0.00
42) 2,4,6-Tribromophenol	10.66	330	378157	82.12	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1831930	62.54	ng	0.00
79) Terphenyl-d14	12.94	244	1547641	61.01	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	338625	10.898	ng	99
71) Phenanthrene	11.38	178	158105	4.524	ng	96
75) Fluoranthene	12.57	202	269215	7.160	ng	96
78) Pyrene	12.79	202	247658	6.308	ng	97
81) Benzo(a)anthracene	13.98	228	132851	4.391	ng	99
83) Chrysene	14.02	228	119425	3.765	ng	96
87) Indeno(1,2,3-cd)pyrene	16.87	276	64871	2.011	ng	95
88) Benzo(b)fluoranthene	15.02	252	158393m	4.836	ng	
90) Benzo(a)pyrene	15.38	252	110873	3.870	ng	98
92) Benzo(g,h,i)perylene	17.31	276	62346	2.404	ng	97

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

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