

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030920\
 Data File : BF119310.D
 Acq On : 10 Mar 2020 4:19
 Operator : CG/JU
 Sample : L1836-01 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-212-VNJ-221

Manual Integrations
 APPROVED

mohammad
 3/10/2020 2:45:04 PM

Quant Time: Mar 10 12:16:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 06 23:38:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	95785	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	349420	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	173939	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	311570	20.00	ng	0.00
76) Chrysene-d12	13.89	240	265991	20.00	ng	0.00
87) Perylene-d12	15.32	264	208427	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	227661	39.72	ng	0.00
7) Phenol-d6	6.40	99	280151	40.19	ng	0.02
23) Nitrobenzene-d5	7.29	82	168750	25.50	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	80469	35.36	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	351343	29.76	ng	0.00
79) Terphenyl-d14	12.82	244	436276	28.24	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.48	163	34970	2.629	ng	98
71) Phenanthrene	11.27	178	51061	3.005	ng	99
75) Fluoranthene	12.46	202	132771	7.361	ng	99
78) Pyrene	12.69	202	117934	5.522	ng	98
81) Benzo(a)anthracene	13.87	228	60279	3.326	ng	97
83) Chrysene	13.91	228	62894	3.483	ng	98
84) Bis(2-ethylhexyl)phthalate	13.87	149	2181675	192.105	ng	# 98
86) Indeno(1,2,3-cd)pyrene	16.74	276	35023	2.003	ng	# 94
88) Benzo(b)fluoranthene	14.91	252	80410m	5.853	ng	
89) Benzo(k)fluoranthene	14.94	252	29790m	2.340	ng	
90) Benzo(a)pyrene	15.26	252	44962	3.626	ng	# 93
92) Benzo(a,h,i)perylene	17.17	276	35577	3.300	ng	# 89

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

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