

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082820\
 Data File : BF121466.D
 Acq On : 28 Aug 2020 18:30
 Operator : JU/CG
 Sample : L3837-16
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 789UMR-TP11-0006-02

Manual Integrations
 APPROVED

mohammad
 8/31/2020 10:27:27 AM

Quant Time: Aug 28 19:13:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 28 09:25:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	299023	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1165383	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	600324	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	1134882	20.00	ng	0.00
76) Chrysene-d12	14.02	240	794120	20.00	ng	0.00
86) Perylene-d12	15.49	264	703420	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1403903	78.77	ng	0.00
7) Phenol-d6	6.48	99	1832667	73.58	ng	-0.01
23) Nitrobenzene-d5	7.42	82	1161850	68.16	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	526362	86.81	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	2074134	55.75	ng	0.00
79) Terphenyl-d14	12.97	244	2117052	56.12	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.61	163	210028	5.093	ng	98
71) Phenanthrene	11.40	178	140036	2.404	ng	95
75) Fluoranthene	12.59	202	322289	5.135	ng	95
78) Pyrene	12.82	202	263102	4.639	ng	95
81) Benzo(a)anthracene	14.01	228	122471	2.526	ng	98
83) Chrysene	14.04	228	140387	2.895	ng	96
88) Benzo(b)fluoranthene	15.06	252	181046m	3.957	ng	
90) Benzo(a)pyrene	15.42	252	112689	2.739	ng	97
92) Benzo(g,h,i)perylene	17.37	276	86341	2.530	ng	98

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

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