

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111819\
 Data File : BF117846.D
 Acq On : 18 Nov 2019 20:16
 Operator : JU
 Sample : K5865-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3-(4-6)

Manual Integrations
 APPROVED

mohammad
 11/19/2019 2:36:24 PM

Quant Time: Nov 19 00:29:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 18:36:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	221913	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	831720	20.00	ng	0.00
39) Acenaphthene-d10	9.97	164	386444	20.00	ng	0.00
64) Phenanthrene-d10	11.46	188	610657	20.00	ng	0.00
76) Chrysene-d12	14.11	240	544226	20.00	ng	0.00
87) Perylene-d12	15.63	264	567003	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	66791	5.33	ng	0.00
7) Phenol-d6	6.53	99	855256	56.56	ng	-0.01
23) Nitrobenzene-d5	7.48	82	941835	67.32	ng	-0.01
42) 2,4,6-Tribromophenol	10.76	330	716	0.18	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	1729508	80.29	ng	0.00
79) Terphenyl-d14	13.05	244	1670011	56.08	ng	0.00
Target Compounds						
10) Phenol	6.55	94	101091	6.434	ng	98
50) Dimethylphthalate	9.67	163	200481	7.398	ng	99
71) Phenanthrene	11.49	178	252888	8.237	ng	95
75) Fluoranthene	12.68	202	322057	4.797	ng	98
78) Pyrene	12.91	202	326161	7.416	ng	98
81) Benzo(a)anthracene	14.10	228	166850	4.639	ng	95
83) Chrysene	14.13	228	165052	4.736	ng	97
84) Bis(2-ethylhexyl)phthalate	14.09	149	54476	2.379	ng	99
86) Indeno(1,2,3-cd)pyrene	17.15	276	63923	2.155	ng	96
88) Benzo(b)fluoranthene	15.17	252	198172m	5.380	ng	
90) Benzo(a)pyrene	15.56	252	131309	4.054	ng	# 92
92) Benzo(a,h,i)perylene	17.62	276	60885	2.192	ng	# 92

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

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