

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111819\
 Data File : BF117845.D
 Acq On : 18 Nov 2019 19:48
 Operator : JU
 Sample : K5865-08
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 3-(2-4)

Manual Integrations
 APPROVED

mohammad
 11/19/2019 2:36:22 PM

Quant Time: Nov 19 00:26:31 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	237584	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	875935	20.00	ng	-0.01
39) Acenaphthene-d10	9.97	164	409997	20.00	ng	0.00
64) Phenanthrene-d10	11.46	188	627961	20.00	ng	0.00
76) Chrysene-d12	14.11	240	514791	20.00	ng	0.00
87) Perylene-d12	15.62	264	561814	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	127649	9.51	ng	0.00
7) Phenol-d6	6.53	99	1040891	64.30	ng	-0.01
23) Nitrobenzene-d5	7.48	82	1029095	69.84	ng	-0.01
42) 2,4,6-Tribromophenol	10.76	330	4200	0.97	ng	-0.01
45) 2-Fluorobiphenyl	9.29	172	1883719	82.43	ng	0.00
79) Terphenyl-d14	13.05	244	1680910	59.68	ng	0.00
Target Compounds						
10) Phenol	6.55	94	113395	6.741	ng	97
15) Benzyl Alcohol	7.05	79	30710	2.457	ng	98
50) Dimethylphthalate	9.67	163	234116	8.143	ng	100
71) Phenanthrene	11.49	178	209531	6.637	ng	96
75) Fluoranthene	12.68	202	283810	2.818	ng	97
78) Pyrene	12.91	202	275978	6.634	ng	99
81) Benzo(a)anthracene	14.10	228	152710	4.489	ng	96
83) Chrysene	14.13	228	138144	4.191	ng	98
84) Bis(2-ethylhexyl)phthalate	14.09	149	50861	2.348	ng	95
86) Indeno(1,2,3-cd)pyrene	17.15	276	63041	2.247	ng	95
88) Benzo(b)fluoranthene	15.17	252	198290m	5.432	ng	
90) Benzo(a)pyrene	15.56	252	121792	3.794	ng	# 91
92) Benzo(a,h,i)perylene	17.62	276	58602	2.130	ng	93

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

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