

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110719\
 Data File : BF117724.D
 Acq On : 8 Nov 2019 00:02
 Operator : JU
 Sample : K5722-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 6-A

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:31:48 PM

Quant Time: Nov 08 03:03:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 17:27:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	296432	20.00	ng	0.00
21) Naphthalene-d8	8.22	136	1147283	20.00	ng	0.00
39) Acenaphthene-d10	9.98	164	596450	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1017647	20.00	ng	0.00
76) Chrysene-d12	14.12	240	695737	20.00	ng	0.00
87) Perylene-d12	15.63	264	843594	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	1461257	89.23	ng	0.00
7) Phenol-d6	6.55	99	1863112	92.27	ng	-0.01
23) Nitrobenzene-d5	7.49	82	1111633	57.20	ng	-0.01
42) 2,4,6-Tribromophenol	10.77	330	498339	85.41	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	2094015	69.27	ng	0.00
79) Terphenyl-d14	13.06	244	1978107	58.53	ng	0.00
Target Compounds						
10) Phenol	6.56	94	65055	2.990	ng	95
50) Dimethylphthalate	9.69	163	244357	6.060	ng	98
71) Phenanthrene	11.49	178	300389	6.158	ng	95
75) Fluoranthene	12.69	202	412005	8.272	ng	96
78) Pyrene	12.92	202	425595	8.380	ng	97
81) Benzo(a)anthracene	14.11	228	187290	4.325	ng	96
83) Chrysene	14.14	228	226928	5.306	ng	96
86) Indeno(1,2,3-cd)pyrene	17.16	276	109073	2.740	ng	94
88) Benzo(b)fluoranthene	15.18	252	285697m	5.377	ng	
90) Benzo(a)pyrene	15.56	252	161700	3.511	ng	96
92) Benzo(g,h,i)perylene	17.62	276	98471	2.460	ng	96

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

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