

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
 Data File : BF118008.D
 Acq On : 26 Nov 2019 15:43
 Operator : JU
 Sample : K5672-01DL 10X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-112219DL

Manual Integrations
 APPROVED

jagrut
 11/27/2019 11:33:34 PM

Quant Time: Nov 26 16:15:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF111319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 05:38:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	154445	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	482589	20.00	ng	0.00
39) Acenaphthene-d10	9.95	164	215785	20.00	ng	0.00
64) Phenanthrene-d10	11.44	188	330921	20.00	ng	0.00
76) Chrysene-d12	14.09	240	248812	20.00	ng	0.00
87) Perylene-d12	15.59	264	209764	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	16174	1.85	ng	0.00
7) Phenol-d6	6.51	99	22384	2.13	ng	-0.01
23) Nitrobenzene-d5	7.45	82	9577	1.18	ng	-0.01
42) 2,4,6-Tribromophenol	10.73	330	3062	1.34	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	21923	1.82	ng	0.00
79) Terphenyl-d14	13.03	244	18389	1.35	ng	0.00
Target Compounds						
58) Fluorene	10.49	166	41441	3.140	ng	Qvalue 97
71) Phenanthrene	11.46	178	223940	13.460	ng	96
72) Anthracene	11.52	178	72098	4.371	ng	96
75) Fluoranthene	12.66	202	191959	6.178	ng	98
78) Pyrene	12.89	202	208331	10.361	ng	98
81) Benzo(a)anthracene	14.08	228	92027	5.597	ng	97
83) Chrysene	14.12	228	75958	4.768	ng	96
86) Indeno(1,2,3-cd)pyrene	17.12	276	28614	2.110	ng	95
88) Benzo(b)fluoranthene	15.15	252	66646m	4.890	ng	
90) Benzo(a)pyrene	15.53	252	50710	4.231	ng	# 93
92) Benzo(a,h,i)perylene	17.57	276	30054	2.925	ng	# 89

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

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