

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102519\
 Data File : BF117542.D
 Acq On : 25 Oct 2019 18:00
 Operator : JU
 Sample : K5512-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CHRT-26706

Manual Integrations
 APPROVED

mohammad
 10/29/2019 3:33:31 PM

Quant Time: Oct 26 01:41:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	34251	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	135703	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	62812	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	85393	20.00	ng	0.00
76) Chrysene-d12	13.89	240	68365	20.00	ng	0.00
87) Perylene-d12	15.33	264	59777	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	244791	106.33	ng	0.00
7) Phenol-d6	6.38	99	334880	108.30	ng	0.00
23) Nitrobenzene-d5	7.29	82	203042	73.87	ng	-0.01
42) 2,4,6-Tribromophenol	10.56	330	49205	90.54	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	342104	76.74	ng	-0.01
79) Terphenyl-d14	12.83	244	260140	62.07	ng	0.00
Target Compounds						
10) Phenol	6.39	94	8209	2.601	ng	# 80
49) Acenaphthylene	9.62	152	12373	2.149	ng	# 95
50) Dimethylphthalate	9.48	163	36732	8.203	ng	99
71) Phenanthrene	11.27	178	35215	7.102	ng	95
72) Anthracene	11.33	178	10804	2.124	ng	98
75) Fluoranthene	12.46	202	25385	5.114	ng	98
78) Pyrene	12.69	202	40965	6.976	ng	98
81) Benzo(a)anthracene	13.88	228	13169	2.856	ng	92
83) Chrysene	13.92	228	13883	3.072	ng	96
88) Benzo(b)fluoranthene	14.93	252	10349m	2.932	ng	
90) Benzo(a)pyrene	15.27	252	10148	3.162	ng	# 81
92) Benzo(a,h,i)perylene	17.18	276	5990	2.281	ng	# 82

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

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