

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102519\
 Data File : BF117535.D
 Acq On : 25 Oct 2019 14:52
 Operator : JU
 Sample : K5502-15
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 178-02-(0-2.3)

Manual Integrations
 APPROVED

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

Quant Time: Oct 26 01:11:39 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	39284	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	153818	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	81800	20.00	ng	-0.01
64) Phenanthrene-d10	11.25	188	127859	20.00	ng	0.00
76) Chrysene-d12	13.89	240	76418	20.00	ng	0.00
87) Perylene-d12	15.32	264	79176	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	248612	94.15	ng	0.00
7) Phenol-d6	6.38	99	338400	95.42	ng	0.00
23) Nitrobenzene-d5	7.29	82	211586	67.91	ng	-0.01
42) 2,4,6-Tribromophenol	10.56	330	67371	95.19	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	333603	57.46	ng	-0.01
79) Terphenyl-d14	12.83	244	256506	54.75	ng	0.00
Target Compounds						
10) Phenol	6.39	94	8828	2.439	ng	95
50) Dimethylphthalate	9.48	163	66430	11.392	ng	100
71) Phenanthrene	11.27	178	58294	7.852	ng	98
75) Fluoranthene	12.46	202	77146	10.380	ng	97
78) Pyrene	12.69	202	64964	9.897	ng	99
81) Benzo(a)anthracene	13.88	228	29853	5.791	ng	99
83) Chrysene	13.92	228	25224	4.993	ng	96
86) Indeno(1,2,3-cd)pyrene	16.74	276	14141	3.749	ng	95
88) Benzo(b)fluoranthene	14.92	252	27680m	5.920	ng	
89) Benzo(k)fluoranthene	14.94	252	9960m	2.078	ng	
90) Benzo(a)pyrene	15.27	252	20571	4.839	ng	# 88
92) Benzo(a,h,i)perylene	17.17	276	13474	3.874	ng	# 89

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

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