

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF103019\
 Data File : BF117630.D
 Acq On : 30 Oct 2019 17:46
 Operator : JU
 Sample : K5542-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 199-48-(0-1.2)

Manual Integrations
 APPROVED

mohammad
 10/31/2019 3:10:53 PM

Quant Time: Oct 31 03:41:55 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	37315	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	149851	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	71249	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	104221	20.00	ng	0.00
76) Chrysene-d12	13.90	240	79500	20.00	ng	0.00
87) Perylene-d12	15.33	264	65105	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	207246	82.63	ng	0.00
7) Phenol-d6	6.39	99	287716	85.41	ng	0.00
23) Nitrobenzene-d5	7.29	82	167995	55.35	ng	-0.01
42) 2,4,6-Tribromophenol	10.56	330	44784	72.64	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	291770	57.70	ng	-0.01
79) Terphenyl-d14	12.83	244	241391	49.53	ng	0.00
Target Compounds						
10) Phenol	6.40	94	7879	2.292	ng	Qvalue # 70
50) Dimethylphthalate	9.48	163	43372	8.539	ng	99
52) Acenaphthene	9.79	154	29108	7.263	ng	99
55) Dibenzofuran	9.97	168	30041	5.174	ng	96
58) Fluorene	10.31	166	46447	9.673	ng	97
71) Phenanthrene	11.27	178	445716	73.651	ng	99
72) Anthracene	11.33	178	145729	23.469	ng	98
73) Carbazole	11.49	167	40784	7.187	ng	99
75) Fluoranthene	12.47	202	495751	81.831	ng	99
78) Pyrene	12.70	202	398332	58.333	ng	100
81) Benzo(a)anthracene	13.89	228	185547	34.600	ng	99
83) Chrysene	13.92	228	149324	28.412	ng	99
86) Indeno(1,2,3-cd)pyrene	16.75	276	53442	13.620	ng	95
88) Benzo(b)fluoranthene	14.93	252	151280m	39.349	ng	
89) Benzo(k)fluoranthene	14.94	252	54478m	13.825	ng	
90) Benzo(a)pyrene	15.27	252	102609	29.355	ng	# 94
91) Dibenzo(a,h)anthracene	16.74	278	11229	3.715	ng	# 75
92) Benzo(g,h,i)perylene	17.18	276	46743	16.345	ng	94

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

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