

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062920\
 Data File : BF120741.D
 Acq On : 29 Jun 2020 21:38
 Operator : JU/CG
 Sample : L3128-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 44-08

Manual Integrations
 APPROVED

mohammad
 6/30/2020 4:14:53 PM

Quant Time: Jun 30 02:24:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 29 15:30:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	50157	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	462865	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	96045	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	170092	20.00	ng	0.01
76) Chrysene-d12	13.96	240	146580	20.00	ng	0.00
86) Perylene-d12	15.41	264	152915	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	253408	86.62	ng	0.08
7) Phenol-d6	6.51	99	326862	89.70	ng	0.08
23) Nitrobenzene-d5	7.36	82	477848	57.71	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	76776	68.76	ng	0.01
45) 2-Fluorobiphenyl	9.15	172	688852	109.73	ng	0.00
79) Terphenyl-d14	12.91	244	488673	73.64	ng	0.01
Target Compounds						
50) Dimethylphthalate	9.55	163	44342	6.410	ng	98
71) Phenanthrene	11.35	178	46336	5.480	ng	96
75) Fluoranthene	12.54	202	100386	10.384	ng	98
78) Pyrene	12.77	202	89414	8.611	ng	99
81) Benzo(a)anthracene	13.96	228	54535	6.193	ng	99
83) Chrysene	13.99	228	51061	5.653	ng	96
84) Bis(2-ethylhexyl)phthalate	13.94	149	45453	8.077	ng	98
87) Indeno(1,2,3-cd)pyrene	16.83	276	35640	3.832	ng	97
88) Benzo(b)fluoranthene	14.99	252	79202m	8.695	ng	
89) Benzo(k)fluoranthene	15.02	252	20950m	2.599	ng	
90) Benzo(a)pyrene	15.34	252	52822	6.480	ng	97
92) Benzo(a,h,i)perylene	17.27	276	34581	4.600	ng	98

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

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