

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072720\
 Data File : BF121036.D
 Acq On : 27 Jul 2020 17:13
 Operator : JU/CG
 Sample : L3383-09
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CRT-009

Manual Integrations
 APPROVED

mohammad
 7/28/2020 11:43:25 AM

Quant Time: Jul 27 18:03:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 16 14:20:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	199029	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	739838	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	410045	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	838263	20.00	ng	0.00
76) Chrysene-d12	13.97	240	681160	20.00	ng	0.00
86) Perylene-d12	15.42	264	598615	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	942014	77.59	ng	0.00
7) Phenol-d6	6.45	99	1137046	72.50	ng	0.00
23) Nitrobenzene-d5	7.37	82	717017	56.08	ng	0.00
42) 2,4,6-Tribromophenol	10.64	330	389906	86.87	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	1430429	52.08	ng	0.00
79) Terphenyl-d14	12.92	244	1535289	49.00	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.56	163	317044	10.717	ng	99
75) Fluoranthene	12.54	202	155500	3.254	ng	99
78) Pyrene	12.77	202	227692	4.728	ng	99
81) Benzo(a)anthracene	13.96	228	82840	2.008	ng	96
83) Chrysene	13.99	228	116809	2.695	ng	96
88) Benzo(b)fluoranthene	14.99	252	151116m	4.176	ng	
90) Benzo(a)pyrene	15.35	252	76801	2.326	ng	99

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

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