

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071520\
 Data File : BP002765.D
 Acq On : 15 Jul 2020 21:19
 Operator : CG/JU
 Sample : L3188-07 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TR-05-071420

Manual Integrations
 APPROVED

mohammad
 7/16/2020 11:07:32 AM

Quant Time: Jul 16 01:51:07 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 13 10:54:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	168526	20.00	ng	0.01
21) Naphthalene-d8	10.33	136	679321	20.00	ng	0.00
39) Acenaphthene-d10	14.22	164	399693	20.00	ng	0.00
64) Phenanthrene-d10	16.97	188	807374	20.00	ng	0.00
76) Chrysene-d12	21.08	240	641589	20.00	ng	0.00
86) Perylene-d12	23.27	264	612673	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	75791	7.59	ng	0.01
7) Phenol-d6	6.77	99	98538	6.91	ng	0.01
23) Nitrobenzene-d5	8.72	82	50333	3.96	ng	0.01
42) 2,4,6-Tribromophenol	15.72	330	27495	6.59	ng	0.01
45) 2-Fluorobiphenyl	12.83	172	138810	5.37	ng	0.00
79) Terphenyl-d14	19.56	244	160898	5.75	ng	0.00
Target Compounds						
75) Fluoranthene	19.00	202	219509	4.334	ng	99
78) Pyrene	19.35	202	194531	4.391	ng	98
81) Benzo(a)anthracene	21.07	228	105938	2.551	ng	# 90
83) Chrysene	21.12	228	93729	2.371	ng	# 93
88) Benzo(b)fluoranthene	22.62	252	132187m	3.475	ng	
90) Benzo(a)pyrene	23.18	252	92289	2.560	ng	# 84

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

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