

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072420\
 Data File : BP002866.D
 Acq On : 24 Jul 2020 22:13
 Operator : CG/JU
 Sample : L3383-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CRT-009

Quant Time: Jul 25 00:32:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 24 15:50:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	31416	20.00	ng	0.00
21) Naphthalene-d8	10.31	136	131639	20.00	ng	0.00
39) Acenaphthene-d10	14.19	164	89710	20.00	ng	0.00
64) Phenanthrene-d10	16.95	188	239314	20.00	ng	0.00
76) Chrysene-d12	21.05	240	365720	20.00	ng	0.00
86) Perylene-d12	23.23	264	450719	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	185606	99.68	ng	0.00
7) Phenol-d6	6.74	99	251159	94.51	ng	0.00
23) Nitrobenzene-d5	8.69	82	141282	57.40	ng	0.00
42) 2,4,6-Tribromophenol	15.70	330	97324	103.93	ng	0.00
45) 2-Fluorobiphenyl	12.80	172	322672	55.64	ng	0.00
79) Terphenyl-d14	19.53	244	761988	47.77	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	13.66	163	66492	9.719	ng	99
75) Fluoranthene	18.97	202	52571	3.501	ng	99
78) Pyrene	19.33	202	83360	3.301	ng	99
83) Chrysene	21.09	228	57659	2.559	ng	96
88) Benzo(b)fluoranthene	22.59	252	107212	3.831	ng	# 97
90) Benzo(a)pyrene	23.14	252	58283	2.198	ng	95

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

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