

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072420\
 Data File : BP002869.D
 Acq On : 24 Jul 2020 23:56
 Operator : CG/JU
 Sample : L3383-01
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CRT-013

Manual Integrations
 APPROVED

mohammad
 7/27/2020 12:22:23 PM

Quant Time: Jul 25 01:06:27 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	37308	20.00	ng	0.00
21) Naphthalene-d8	10.30	136	163211	20.00	ng	0.00
39) Acenaphthene-d10	14.19	164	114075	20.00	ng	0.00
64) Phenanthrene-d10	16.95	188	285804	20.00	ng	0.00
76) Chrysene-d12	21.05	240	405386	20.00	ng	0.00
86) Perylene-d12	23.23	264	491638	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	250345	113.21	ng	0.00
7) Phenol-d6	6.74	99	357468	113.27	ng	0.00
23) Nitrobenzene-d5	8.68	82	205881	67.47	ng	0.00
42) 2,4,6-Tribromophenol	15.69	330	142606	119.76	ng	0.00
45) 2-Fluorobiphenyl	12.80	172	491722	66.68	ng	0.00
79) Terphenyl-d14	19.53	244	1007553	56.98	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.65	163	96813	11.128	ng	99
70) Pentachlorophenol	16.63	266	4809	10.191	ng	95
71) Phenanthrene	16.99	178	60497	3.909	ng	99
75) Fluoranthene	18.97	202	276496	15.420	ng	99
78) Pyrene	19.33	202	321847	11.499	ng	99
81) Benzo(a)anthracene	21.04	228	158115	6.025	ng	95
83) Chrysene	21.09	228	165588	6.629	ng	94
87) Indeno(1,2,3-cd)pyrene	25.42	276	79996	2.251	ng	97
88) Benzo(b)fluoranthene	22.59	252	243041	7.961	ng	98
89) Benzo(k)fluoranthene	22.62	252	80109m	2.664	ng	
90) Benzo(a)pyrene	23.14	252	128822	4.453	ng	# 94
92) Benzo(a,h,i)perylene	26.09	276	73975	2.546	ng	98

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

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