

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP072220\
 Data File : BP002834.D
 Acq On : 22 Jul 2020 20:23
 Operator : CG/JU
 Sample : L3360-01RX
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SP-1ARX

Manual Integrations
 APPROVED

Sohil
 7/24/2020 12:11:06 PM

Quant Time: Jul 23 01:47:54 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	7.55	152	176449	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	693093	20.00	ng	-0.01
39) Acenaphthene-d10	14.20	164	383786	20.00	ng	-0.01
64) Phenanthrene-d10	16.96	188	700600	20.00	ng	0.00
76) Chrysene-d12	21.07	240	529894	20.00	ng	0.00
86) Perylene-d12	23.26	264	530414	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	265561	25.39	ng	0.00
7) Phenol-d6	6.76	99	792280	53.08	ng	0.00
23) Nitrobenzene-d5	8.69	82	673077	51.94	ng	-0.01
42) 2,4,6-Tribromophenol	15.71	330	66092	16.50	ng	0.00
45) 2-Fluorobiphenyl	12.82	172	1362503	54.92	ng	0.00
79) Terphenyl-d14	19.55	244	1257292	54.40	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.66	163	211766	7.235	ng	99
71) Phenanthrene	17.00	178	227728	6.003	ng	100
75) Fluoranthene	18.99	202	602216	13.701	ng	98
78) Pyrene	19.35	202	529363	14.469	ng	100
81) Benzo(a)anthracene	21.06	228	263841	7.692	ng	96
83) Chrysene	21.11	228	249590	7.644	ng	96
87) Indeno(1,2,3-cd)pyrene	25.47	276	125101	3.263	ng	97
88) Benzo(b)fluoranthene	22.62	252	283669	8.613	ng	# 94
89) Benzo(k)fluoranthene	22.65	252	96478m	2.974	ng	
90) Benzo(a)pyrene	23.17	252	200208	6.415	ng	# 96
92) Benzo(g,h,i)perylene	26.15	276	123878	3.952	ng	97

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

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