

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP073020\
 Data File : BP002886.D
 Acq On : 30 Jul 2020 20:42
 Operator : CG/JU
 Sample : L3491-12
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-3-S

Manual Integrations
 APPROVED

Sohil
 7/31/2020 12:11:00 PM

Quant Time: Jul 31 01:38:21 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 30 15:30:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	550455	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	2041616	20.00	ng	0.00
39) Acenaphthene-d10	14.35	164	1186299	20.00	ng	0.00
64) Phenanthrene-d10	17.10	188	2553486	20.00	ng	0.00
76) Chrysene-d12	21.20	240	2724513	20.00	ng	0.00
86) Perylene-d12	23.46	264	2778217	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	2483044	69.04	ng	0.00
7) Phenol-d6	6.89	99	3113467	62.35	ng	0.00
23) Nitrobenzene-d5	8.86	82	1867778	49.55	ng	0.00
42) 2,4,6-Tribromophenol	15.85	330	1025878	69.58	ng	0.00
45) 2-Fluorobiphenyl	12.97	172	3921124	45.16	ng	0.00
79) Terphenyl-d14	19.68	244	5422831	39.85	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.81	163	579026	6.241	ng	99
71) Phenanthrene	17.15	178	381946	2.551	ng	98
75) Fluoranthene	19.12	202	1494446	8.663	ng	99
78) Pyrene	19.47	202	1403465	7.369	ng	98
81) Benzo(a)anthracene	21.18	228	1020813	5.501	ng	99
83) Chrysene	21.23	228	820025	4.684	ng	96
87) Indeno(1,2,3-cd)pyrene	25.77	276	585199	2.630	ng	96
88) Benzo(b)fluoranthene	22.78	252	1180117m	6.019	ng	
89) Benzo(k)fluoranthene	22.82	252	434811m	2.346	ng	
90) Benzo(a)pyrene	23.36	252	904510	5.106	ng	99
92) Benzo(g,h,i)perylene	26.47	276	606639	3.313	ng	97

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

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