

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112021\
 Data File : BP008058.D
 Acq On : 20 Nov 2021 15:27
 Operator : CG/JU
 Sample : M4724-28
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 T-132-2

Quant Time: Nov 22 05:31:41 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 20 00:21:11 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	121430	20.000	ng	0.00
21) Naphthalene-d8	10.616	136	466552	20.000	ng	-0.01
39) Acenaphthene-d10	14.463	164	307377	20.000	ng	0.00
64) Phenanthrene-d10	17.216	188	640603	20.000	ng	-0.01
76) Chrysene-d12	21.310	240	652115	20.000	ng	-0.01
86) Perylene-d12	23.704	264	740040	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.411	112	846921	119.213	ng	0.00
7) Phenol-d6	6.999	99	1120372	113.769	ng	-0.01
23) Nitrobenzene-d5	8.981	82	745699	72.508	ng	-0.01
42) 2,4,6-Tribromophenol	15.957	330	389224	100.059	ng	-0.01
45) 2-Fluorobiphenyl	13.087	172	1521388	77.503	ng	0.00
79) Terphenyl-d14	19.786	244	2334917	70.396	ng	-0.01
Target Compounds						
						Qvalue
74) Di-n-butylphthalate	18.181	149	101059	2.672	ng	96
87) Indeno(1,2,3-cd)pyrene	26.198	276	149840	2.918	ng	99
88) Benzo(b)fluoranthene	22.974	252	150013	3.242	ng	# 98
89) Benzo(k)fluoranthene	23.021	252	157681	3.425	ng	# 95
90) Benzo(a)pyrene	23.598	252	154945	3.996	ng	# 97
91) Dibenzo(a,h)anthracene	26.210	278	103499	2.421	ng	# 94
92) Benzo(g,h,i)perylene	26.951	276	103441	2.470	ng	98

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

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